25

30

What is claimed is:

1. A compound of the formula (I):

(I)

or a stereoisomer or a pharmaceutically acceptable salt form thereof, wherein:

b is a single bond wherein the bridging hydrogens are either cis or trans:

15 X is a bond, -CH₂-, -O-, -S-, -S(=O)-, -S(=O) $_2$ -, -NR¹⁰-,

 $-CH_2CH_2-$, $-OCH_2-$, $-SCH_2-$, $-S(=O)CH_2-$, $-S(=O)_2CH_2-$,

-CH₂O-, -CH₂S-, -CH₂S(=O)-, -CH₂S(=O)₂-, -NR¹⁰CH₂-,

-CH $_2$ NR 10 -, -NHC(=O)-, or -C(=O)NH-;

20 R^1 is selected from

Н,

 $C(=0)R^{2}$.

 $C(=0)OR^2$.

 C_{1-8} alkyl,

- - - - - -

 C_{2-8} alkenyl,

C₂₋₈ alkynyl, C₃₋₇ cycloalkyl,

C1-6 alkyl substituted with Z,

 C_{2-6} alkenyl substituted with Z,

C2-6 alkynyl substituted with Z,

C₃₋₆ cycloalkyl substituted with Z,

aryl substituted with Z.

5-6 membered heterocyclic ring system containing at least one heteroatom selected from the group

```
consisting of N, O, and S, said heterocyclic ring
                 system substituted with Z;
            C<sub>1-3</sub> alkyl substituted with Y,
            C2-3 alkenyl substituted with Y,
 5
            C_{2-3} alkynyl substituted with Y,
            C<sub>1-6</sub> alkyl substituted with 0-2 R<sup>2</sup>.
            C2-6 alkenyl substituted with 0-2 R2,
            C_{2-6} alkynyl substituted with 0-2 R^2.
            aryl substituted with 0-2 R2, and
10
            5-6 membered heterocyclic ring system containing at
                 least one heteroatom selected from the group
                 consisting of N, O, and S, said heterocyclic ring
                 system substituted with 0-2 R2;
15
     Y is selected from
            C<sub>3-6</sub> cycloalkyl substituted with Z.
            aryl substituted with Z,
            5-6 membered heterocyclic ring system containing at
                 least one heteroatom selected from the group
                consisting of N, O, and S, said heterocyclic ring
20
                system substituted with Z;
            C_{3-6} cycloalkyl substituted with -(C_{1-3} alkyl)-Z,
            aryl substituted with -(C_{1-3} \text{ alkyl}) - Z, and
            5-6 membered heterocyclic ring system containing at
25
                least one heteroatom selected from the group
                consisting of N, O, and S, said heterocyclic ring
                system substituted with -(C_{1-3} \text{ alkyl}) - Z;
     Z is selected from H.
30
          -CH (OH) R2,
          -C(ethylenedioxy)R2.
          -OR2,
          -SR2,
          -NR^2R^3.
35
          -C(0)R2.
          -C(O)NR2R3.
          -NR^3C(0)R^2
```

```
-C(O)OR2,
            -OC(O)R2,
            -CH(=NR^4)NR^2R^3,
            -NHC (=NR^4) NR^2R^3.
            -S(O)R2,
  5
            -S(0)_2R^2,
            -S(O)_2NR^2R^3, and -NR^3S(O)_2R^2;
      \mathbb{R}^2, at each occurrence, is independently selected from
10
            halo.
            C_{1-3} haloalkyl,
           C_{1-4} alkyl,
           C_{2-4} alkenyl,
           C2-4 alkynyl,
15
           C3-6 Cycloalkyl,
           aryl substituted with 0-5 R42;
           C_{3-10} carbocyclic residue substituted with 0-3 R^{41}, and
           5-10 membered heterocyclic ring system containing from
                 1-4 heteroatoms selected from the group
                 consisting of N, O, and S substituted with 0-3
20
                 R41:
     \mathbb{R}^3, at each occurrence, is independently selected from
            H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, and
25
           C_{1-4} alkoxy;
     alternatively, {\bf R}^2 and {\bf R}^3 join to form a 5- or 6-membered
           ring optionally substituted with -O- or -N(\mathbb{R}^4)-;
30
     \mathbb{R}^4, at each occurrence, is independently selected from H
           and C_{1-4} alkyl;
     R6a is H or C1-4 alkyl;
35
     R6b is H:
```

13

```
alternatively, R^{6a} and R^{6b} are taken together to form =0 or
             =S:
      \ensuremath{\mbox{R}^{7}} and \ensuremath{\mbox{R}^{9}}, at each occurrence, are independently selected
  5
             from
             H, halo, -CF3, -OCF3, -OH, -CN, -NO2, -NR46R47,
             C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl,
                    C_{1-8} alkoxy, (C_{1-4} haloalkyl)oxy,
             C_{3-10} cycloalkyl substituted with 0-2 R^{33},
10
             C_{1-4} alkyl substituted with 0-2 R^{11}.
             C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>,
             aryl substituted with 0-5 R33.
             5-10 membered heterocyclic ring system containing from
                    1-4 heteroatoms selected from the group
15
                   consisting of N, O, and S substituted with 0-3
                   R31;
            OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13},
            NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12}.
20
            CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, S(O)_2R^{12},
            S(0) NR^{12}R^{13}, S(0)_2NR^{12}R^{13}, NR^{14}S(0)R^{12}, NR^{14}S(0)_2R^{12}.
            NR^{12}C(0)R^{15}, NR^{12}C(0)OR^{15}, NR^{12}S(0)_2R^{15}, and
            NR12C(0)NHR15;
25
      R8 is selected from
            H, halo, -CF3, -OCF3, -OH, -CN, -NO2,
            C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, C<sub>1-4</sub> haloalkyl,
                   C<sub>1-8</sub> alkoxy, (C<sub>1-4</sub> haloalkyl) oxy,
            C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>,
30
            C_{1-4} alkyl substituted with 0-2 R^{11},
            C2-4 alkenyl substituted with 0-2 R11.
            C2-4 alkynyl substituted with 0-1 R11,
            C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>,
            aryl substituted with 0-5 R33,
```

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group

```
R31;
               OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13}.
               NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12},
     5
               CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, S(O)_2R^{12}.
               S(O)NR^{12}R^{13}, S(O)_2NR^{12}R^{13}, NR^{14}S(O)R^{12}, NR^{14}S(O)_2R^{12},
               NR^{12}C(O)R^{15}, NR^{12}C(O)OR^{15}, NR^{12}S(O)_2R^{15}, and
               NR12C(O)NHR15:
    10
         R10 is selected from H.
               C1-4 alkyl substituted with 0-2 R10A,
               C2-4 alkenyl substituted with 0-2 R10A.
               C_{2-4} alkynyl substituted with 0-1 R^{10A}, and
   15
               C_{1-4} alkoxy;
         R10A is selected from
               C<sub>1-4</sub> alkoxy,
               C_{3-6} carbocyclic residue substituted with 0-3 R^{33},
1 20
               phenyl substituted with 0-3 R33, and
               5-6 membered heterocyclic ring system containing 1, 2,
                    or 3 heteroatoms selected from the group
                    consisting of N, O, and S; substituted with 0-2
                     R44;
   25
        R11 is selected from
              H, halo, -CF3, -CN, -NO2,
              C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl,
                    C1-8 alkoxy, C3-10 cycloalkyl,
   30
              C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
              aryl substituted with 0-5 R33,
              5-10 membered heterocyclic ring system containing from
                    1-4 heteroatoms selected from the group
                    consisting of N, O, and S substituted with 0-3
   35
                    R31.
```

i sh

13

8 je sk

F.J

So the

0

į, i

consisting of N, O, and S substituted with 0-3 $\,$

10

15

30

R31;

```
OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13},
NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12},
CH (=NR^{14}) NR^{12}R^{13}, NHC (=NR^{14}) NR^{12}R^{13}, S(0) R^{12}, S(0) 2R^{12},
S(O)NR^{12}R^{13}, S(O)_2NR^{12}R^{13}, NR^{14}S(O)R^{12}, NR^{14}S(O)_2R^{12},
NR^{12}C(0)R^{15}, NR^{12}C(0)OR^{15}, NR^{12}S(0)_2R^{15}, and
NR12C (O) NHR15:
```

- \mathbb{R}^{12} , at each occurrence, is independently selected from C₁₋₄ alkyl substituted with 0-1 R^{12a}, C2-4 alkenyl substituted with 0-1 R12a, C2-4 alkynyl substituted with 0-1 R12a. C₃₋₆ cycloalkyl substituted with 0-3 R³³. aryl substituted with 0-5 R33; C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3
- R^{12a} , at each occurrence, is independently selected from phenyl substituted with 0-5 R33; C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group 25 consisting of N, O, and S substituted with 0-3R31.
 - R¹³, at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
 - alternatively, ${\bf R}^{12}$ and ${\bf R}^{13}$ join to form a 5- or 6-membered ring optionally substituted with -O- or -N(R14)-;
- alternatively, \mathbb{R}^{12} and \mathbb{R}^{13} when attached to N may be 35 combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group consisting of N,

30

O, and S, wherein said bicyclic heterocyclic ring system is unsaturated or partially saturated, wherein said bicyclic heterocyclic ring system is substituted with 0-3 R^{16} :

- $R^{14}, \mbox{ at each occurrence, is independently selected from H and <math display="inline">C_{1-4}$ alkyl;
- R^{15} , at each occurrence, is independently selected from 10 H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
 - R^{16} , at each occurrence, is independently selected from H, OH, halo, CN, NO₂, CF_3 , SO_2R^{45} , $NR^{46}R^{47}$, -C(=0)H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl, C_{1-3} haloalkyl-oxy-, C_{1-3} alkyloxy-, and =0;
 - $R^{31},$ at each occurrence, is independently selected from H, OH, halo, $CF_3,\ SO_2R^{45},\ NR^{46}R^{47},\ C_{1-4}\ alkyl,\ and\ =0;$
- 20 R³³, at each occurrence, is independently selected from H, OH, halo, CN, NO₂, CF₃, SO₂R⁴⁵, NR⁴⁶R⁴⁷, -C(=0)H, =0, phenyl, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyl-oxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(=0)-, C₁₋₄ alkyl-C(=0)NH-, C₁₋₄ alkyl-OC(=0)-, C₁₋₄ alkyl-C(=0)NH-, C₁₋₄ alkyl-OC(=0)-,
- C₁₋₄ alkyl-C(=0)O-, C₃₋₆ cycloalkyl-oxy-,
 C₃₋₆ cycloalkylmethyl-oxy-;
 C₁₋₆ alkyl substituted with OH, methoxy, ethoxy,
 propoxy, butoxy, -SO₂R⁴⁵, -NR⁴⁶R⁴⁷, NR⁴⁶R⁴⁷C(=O)-, or
- O (C₁₋₄ alky1)CO₂-; and C₂₋₆ alkenyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, -SO₂R⁴⁵, -NR⁴⁶R⁴⁷, NR⁴⁶R⁴⁷C(=O)-, or (C₁₋₄ alky1)CO₂-;
- 35 R^{41} , at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO₂R⁴⁵, NR⁴⁶R⁴⁷, NO₂, CN, =O; C2₋₈ alkenyl, C2₋₈ alkynyl, C1₋₄ alkoxy, C1₋₄ haloalkyl

```
C<sub>1-4</sub> alkyl substituted with 0-1 R<sup>43</sup>.
              aryl substituted with 0-3 R42, and
             5-10 membered heterocyclic ring system containing from
                    1-4 heteroatoms selected from the group
  5
                    consisting of N, O, and S substituted with 0-3
                    R^{44};
      {\bf R}^{42}, at each occurrence, is independently selected from
             H, CF<sub>3</sub>, halo, OH, CO<sub>2</sub>H, SO<sub>2</sub>R<sup>45</sup>, SOR<sup>45</sup>, SR<sup>45</sup>, NR<sup>46</sup>SO<sub>2</sub>R<sup>45</sup>.
                   NR^{46}COR^{45}, NR^{46}R^{47}, NO_2, CN, CH (=NH) NH_2,
 10
                   NHC (=NH) NH2.
             \text{C}_{2\text{-}6} alkenyl, \text{C}_{2\text{-}6} alkynyl, \text{C}_{1\text{-}4} alkoxy, \text{C}_{1\text{-}4} haloalkyl,
                   C3-6 cycloalkyl,
             C_{1-4} alkyl substituted with 0-1 R^{43},
15
             aryl substituted with 0-3 R^{44}, and
             5-10 membered heterocyclic ring system containing from
                   1-4 heteroatoms selected from the group
                   consisting of N, O, and S substituted with 0-3
                   R44;
20
      R^{43} is C_{3-6} cycloalkyl or aryl substituted with 0-3 R^{44};
      {\bf R}^{44}, at each occurrence, is independently selected from H,
            halo, -OH, NR^{46}R^{47}, CO_2H, SO_2R^{45}, -CF_3, -OCF_3, -CN, -
25
            NO_2, C_{1-4} alkyl, and C_{1-4} alkoxy;
      R^{45} is C_{1-4} alkyl;
      {\rm R}^{46}, at each occurrence, is independently selected from H
30
            and C_{1-4} alkyl;
     {\bf R}^{47}, at each occurrence, is independently selected from H,
            C_{1-4} alkyl, -C(=0) NH(C_{1-4} alkyl), -SO_2(C_{1-4} alkyl),
            -C(=0)0(C_{1-4} alkyl), -C(=0)(C_{1-4} alkyl), and -C(=0)H;
35
     n is 1 or 2:
     m is 1 or 2: and
```

```
n plus m is 2, 3, or 4;
```

provided when n is 1, m is 2, and R^7 , R^8 , and R^9 are independently selected from H, halogen, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkylthio or trifluoromethyl; then X is not a bond.

2. A compound of Claim 1 wherein:

```
10
     X is a bond, -CH_2-, -O-, -S-, -S(=O)-, -S(=O)_2-, -NR^{10}-,
            -CH_2CH_2-, -OCH_2-, -SCH_2-, -CH_2O-, -CH_2S-, -NR^{10}CH_2-, or
            -CH2NR10-:
     R1 is selected from
15
             н.
             C(=0)R^{2}.
             C(=0)OR^2,
             C<sub>1-8</sub> alkyl,
             C2-8 alkenyl,
20
            C2-8 alkynyl,
            C3-7 cycloalkyl,
            C<sub>1-6</sub> alkyl substituted with 0-2 R<sup>2</sup>,
             C_{2-6} alkenyl substituted with 0-2 R^2,
             C2-6 alkynyl substituted with 0-2 R2,
```

aryl substituted with 0-2 R², and
5-6 membered heterocyclic ring system containing at
least one heteroatom selected from the group
consisting of N, O, and S, said heterocyclic ring
system substituted with 0-2 R²;

30

 $\rm R^2,$ at each occurrence, is independently selected from F, Cl, CH₂F, CHF₂, CF₃, C₁₋₄ alkyl, C₂₋₄ alkenyl, C₂₋₄ alkynyl,

35 C₂₋₄ alkynyl,
 C₃₋₆ cycloalkyl,
 phenyl substituted with 0-5 R⁴²;

```
C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>41</sup>, and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R<sup>41</sup>;
```

 R^{6a} is H or C_{1-4} alkyl;

R6b is H;

5

10

15

20

25

30

alternatively, R^{6a} and R^{6b} are taken together to form =0 or =S;

 $\rm R^{7}$ and $\rm R^{9}$, at each occurrence, are independently selected from H, halo, -CF₃, -OCF₃, -OH, -CN, -NO₂, -NR⁴⁶R⁴⁷, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₁₋₄ haloalkyl, C₁₋₈ alkoxy, (C₁₋₄ haloalkyl)oxy, C₃₋₁₀ cycloalkyl substituted with 0-2 R³³, C₁₋₄ alkyl substituted with 0-2 R¹¹,

 C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , arvl substituted with 0-5 R^{33} .

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 \mathbb{R}^{31} :

OR¹², SR¹², NR¹²R¹³, C(O)H, C(O)R¹², C(O)NR¹²R¹³, NR¹⁴C(O)R¹², C(O)OR¹², OC(O)R¹², OC(O)OR¹², CH(=NR¹⁴)NR¹²R¹³, NHC(=NR¹⁴)NR¹²R¹³, S(O)R¹², S(O)R¹², S(O)R¹², S(O)R¹²R¹³, NR¹⁴S(O)R¹², NR¹⁴S(O)R¹², NR¹²C(O)R¹⁵, NR¹²C(O)OR¹⁵, NR¹²S(O)R¹⁵, NR¹²S(O

35 R⁸ is selected from
H, halo, -CF₃, -OCF₃, -OH, -CN, -NO₂,

```
C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl,
                    C_{1-8} alkoxy, (C_{1-4} haloalkyl)oxy,
             C3-10 cycloalkyl substituted with 0-2 R33,
             C1-4 alkyl substituted with 0-2 R11.
  5
             C2-4 alkenyl substituted with 0-2 R11.
             C2-4 alkynyl substituted with 0-1 R11.
             C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>.
             aryl substituted with 0-5 R33.
             5-10 membered heterocyclic ring system containing from
 10
                    1-4 heteroatoms selected from the group
                   consisting of N, O, and S substituted with 0-3
                   R31;
             OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13}.
15
             NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12},
             CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, S(O)_{2}R^{12}.
             S(O)NR^{12}R^{13}, S(O)_2NR^{12}R^{13}, NR^{14}S(O)R^{12}, NR^{14}S(O)_2R^{12},
             NR^{12}C(0)R^{15}, NR^{12}C(0)OR^{15}, NR^{12}S(0)_2R^{15}, and
             NR12C(O)NHR15:
20
      \mbox{R}^{10} is selected from H, \mbox{C}_{1\text{--}4} alkyl, \mbox{C}_{2\text{--}4} alkenyl, \mbox{C}_{2\text{--}4}
             alkynyl, and C1-4 alkoxy;
      R11 is selected from
            H, halo, -CF3, -CN, -NO2,
```

 C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl, C₁₋₈ alkoxy, C₃₋₁₀ cycloalkyl,

 C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , aryl substituted with 0-5 R33,

- 30 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R31;
- 35 OR^{12} , SR^{12} , $NR^{12}R^{13}$, C(O)H, $C(O)R^{12}$, $C(O)NR^{12}R^{13}$. $NR^{14}C(O)R^{12}$, $C(O)OR^{12}$, $OC(O)R^{12}$, $OC(O)OR^{12}$, $CH(=NR^{14})NR^{12}R^{13}$, $NHC(=NR^{14})NR^{12}R^{13}$, $S(0)R^{12}$, $S(0)_{2}R^{12}$.

```
\begin{split} &S\text{(O)} \, NR^{12}R^{13}, \, \, S\text{(O)} \, {}_{2}NR^{12}R^{13}, \, \, NR^{14}S\text{(O)} \, R^{12}, \, \, NR^{14}S\text{(O)} \, {}_{2}R^{12}, \\ &NR^{12}C\text{(O)} \, R^{15}, \, \, NR^{12}C\text{(O)} \, OR^{15}, \, \, NR^{12}S\text{(O)} \, {}_{2}R^{15}, \, \, \text{and} \\ &NR^{12}C\text{(O)} \, NHR^{15}; \end{split}
```

5 R¹², at each occurrence, is independently selected from C₁₋₄ alkyl substituted with 0-1 R^{12a}, C₂₋₄ alkenyl substituted with 0-1 R^{12a}, C₂₋₄ alkynyl substituted with 0-1 R^{12a}, C₃₋₆ cycloalkyl substituted with 0-3 R³³, aryl substituted with 0-5 R³³; C₃₋₁₀ carbocyclic residue substituted with 0-3 R³³, and 5-10 membered heterocyclic ring system containing from

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 $\rm R^{31}$;

 R^{12a} , at each occurrence, is independently selected from phenyl substituted with 0-5 R^{33} ; C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R^{31} ;

25 R^{13} , at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;

alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or -N(R^{14})-;

alternatively, R¹² and R¹³ when attached to N may be
combined to form a 9- or 10-membered bicyclic
heterocyclic ring system containing from 1-3
heteroatoms selected from the group consisting of N,
0, and S, wherein said bicyclic heterocyclic ring
system is unsaturated or partially saturated, wherein

20

said bicyclic heterocyclic ring system is substituted with 0-3 $\ensuremath{\mathrm{R}^{16}}\xspace$;

- R^{14} , at each occurrence, is independently selected from H and C_{1-4} alkyl;
 - $R^{15},$ at each occurrence, is independently selected from $H, \quad C_{1\text{-}4} \text{ alkyl}, \quad C_{2\text{-}4} \text{ alkenyl}, \text{ and } \quad C_{2\text{-}4} \text{ alkynyl};$
- 10 R^{16} , at each occurrence, is independently selected from H, OH, halo, CN, NO₂, CF₃, SO₂R⁴⁵, NR⁴⁶R⁴⁷, -C(=0)H, C₁₋₄ alkyl, C₂₋₄ alkenyl, C₂₋₄ alkynyl, C₁₋₄ haloalkyl, C₁₋₃ haloalkyl-oxy-, C₁₋₃ alkyloxy-, and =0;
 - $\rm R^{31},$ at each occurrence, is independently selected from H, OH, halo, CF3, $\rm SO_2R^{45},\ NR^{46}R^{47},\ C_{1-4}$ alkyl, and =0;
 - $$\begin{split} & R^{33}, \text{ at each occurrence, is independently selected from } \\ & H, \text{ OH, halo, CN, NO}_2, \text{ CF}_3, \text{ SO}_2\text{R}^{45}, \text{ NR}^{46}\text{R}^{47}, \text{ -C(=0)H,} \\ & = \text{O, phenyl, C}_{1-6} \text{ alkyl, C}_{2-6} \text{ alkenyl, C}_{2-6} \text{ alkynyl,} \\ & \text{C}_{3-6} \text{ cycloalkyl, C}_{1-4} \text{ haloalkyl, C}_{1-4} \text{ haloalkyl-oxy-,} \\ & \text{C}_{1-4} \text{ alkyloxy-, C}_{1-4} \text{ alkylthio-, C}_{1-4} \text{ alkyl-C(=0)-,} \\ & \text{C}_{1-4} \text{ alkyl-C(=0)NH-, C}_{1-4} \text{ alkyl-OC(=0)-,} \\ & \text{C}_{1-4} \text{ alkyl-C(=0)0-, C}_{3-6} \text{ cycloalkyl-oxy-,} \end{split}$$
- 25 C_{3-6} cycloalkylmethyl-oxy-; C_{1-6} alkyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, $-SO_2R^{45}$, $-NR^{46}R^{47}$, $NR^{46}R^{47}C$ (=0)-, or $(C_{1-4}$ alkyl) CO_2 -; and
- C₂₋₆ alkenyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, $-SO_2R^{45}$, $-NR^{46}R^{47}$, $NR^{46}R^{47}C(=0)$ -, or (C₁₋₄ alkyl)CO₂-;
- R⁴¹, at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO₂R⁴⁵, NR⁴⁶R⁴⁷, NO₂, CN;

 35 C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₁₋₄ alkoxy, C₁₋₄ haloalkyl C₁₋₄ alkyl substituted with 0-1 R⁴³,

 aryl substituted with 0-3 R⁴², and

```
The state of the s
```

```
5-10 membered heterocyclic ring system containing from
                 1-4 heteroatoms selected from the group
                 consisting of N, O, and S substituted with 0-3
                 R^{44}:
 5
     R42, at each occurrence, is independently selected from
           H, CF<sub>3</sub>, halo, OH, CO<sub>2</sub>H, SO<sub>2</sub>R<sup>45</sup>, NR<sup>46</sup>R<sup>47</sup>, NO<sub>2</sub>, CN,
                 CH (=NH) NH2, NHC (=NH) NH2,
           C2-6 alkenyl, C2-6 alkynyl, C1-4 alkoxy, C1-4 haloalkyl,
10
                 C3-6 cycloalkyl,
           C_{1-4} alkyl substituted with 0-1 R^{43},
           aryl substituted with 0-3 R44, and
           5-10 membered heterocyclic ring system containing from
                 1-4 heteroatoms selected from the group
                 consisting of N. O. and S substituted with 0-3
15
                 R44;
     R43 is C3-6 cycloalkyl or aryl substituted with 0-3 R44;
     R44, at each occurrence, is independently selected from H,
20
           halo, -OH, NR^{46}R^{47}, CO_2H, SO_2R^{45}, -CF<sub>3</sub>, -OCF<sub>3</sub>, -CN, -
           NO_2, C_{1-4} alkyl, and C_{1-4} alkoxy;
     R^{45} is C_{1-4} alkyl;
25
     R46, at each occurrence, is independently selected from H
           and C1-4 alkyl;
     R47, at each occurrence, is independently selected from H
30
           and C_{1-4} alkyl;
     n is 1 or 2:
     m is 1 or 2; and
     n plus m is 2, 3, or 4;
35
     provided when n is 1, m is 2, and \mathbb{R}^7, \mathbb{R}^8, and \mathbb{R}^9 are
     independently selected from H, halogen, C1-4 alkyl, C1-4
```

R6b is H;

alkoxy, C1-4 alkylthio or trifluoromethyl; then X is not a bond.

```
3. A compound of Claim 2 wherein:
                 5
                               X is a bond, -CH_2-, -O-, -S-, -CH_2CH_2-, -OCH_2-, -SCH_2-,
                                                    -CH<sub>2</sub>O-, or -CH<sub>2</sub>S-;
                          R1 is selected from
             1.0
                                                        н,
                                                        C(=0)R^{2}
                                                        C(=0)OR^2
                                                        C_{1-6} alkyl,
                                                    C_{2-6} alkenyl,
                                                       C_{2-6} alkynyl,
                                                       C3-6 cycloalkyl,
The second of th
                                                      C1-4 alkyl substituted with 0-2 R2,
                                                        C_{2-4} alkenyl substituted with 0-2 \mathbb{R}^2, and
                                                        C_{2-4} alkynyl substituted with 0-2 R^2;
                               R2, at each occurrence, is independently selected from
                                                    C_{1-4} alkyl,
                                                    C2-4 alkenyl,
              25
                                                    C2-4 alkynyl,
                                                    C3-6 cycloalkyl,
                                                    phenvl substituted with 0-5 R42;
                                                    C_{3-10} carbocyclic residue substituted with 0-3 \mathbb{R}^{41}, and
                                                    5-10 membered heterocyclic ring system containing from
                                                                         1-4 heteroatoms selected from the group
             30
                                                                         consisting of N, O, and S substituted with 0-3
                                                                         R41;
                               R6a is H or C1-4 alkyl;
```

```
=S;
      R^7 and R^9, at each occurrence, are independently selected
  5
             from
            H, halo, -CF3, -OCF3, -OH, -CN, -NO2, -NR46R47,
             C<sub>1-6</sub> alkyl, C<sub>2-6</sub> alkenyl, C<sub>2-6</sub> alkynyl, C<sub>1-6</sub> haloalkyl,
                   C_{1-6} alkoxy, (C_{1-4} haloalkyl)oxy,
            C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>,
10
            C<sub>1-4</sub> alkyl substituted with 0-2 R<sup>11</sup>,
            C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>.
            aryl substituted with 0-5 R33,
            5-10 membered heterocyclic ring system containing from
                   1-4 heteroatoms selected from the group
15
                   consisting of N, O, and S substituted with 0-3
                   \mathbb{R}^{31}:
            OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13},
                   NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12}.
20
                   CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(0)R^{12}.
                   S(O)_2R^{12}, S(O)NR^{12}R^{13}, S(O)_2NR^{12}R^{13}, NR^{14}S(O)R^{12},
                   and NR14S(0) 2R12:
      R8 is selected from
25
            H, halo, -CF3, -OCF3, -OH, -CN, -NO2,
            C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} haloalkyl,
                  C<sub>1-6</sub> alkoxy, (C<sub>1-4</sub> haloalkyl)oxy,
            C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>,
            C1-4 alkyl substituted with 0-2 R11,
30
            C2-4 alkenyl substituted with 0-2 R11.
            C2-4 alkynyl substituted with 0-1 R11,
            C3-10 carbocyclic residue substituted with 0-3 R33.
            aryl substituted with 0-5 R33.
            5-10 membered heterocyclic ring system containing from
35
                   1-4 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
```

alternatively, R6a and R6b are taken together to form =0 or

R31:

```
OR12, SR12, NR12R13, C(O)H, C(O)R12, C(O)NR12R13,
            NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12},
            CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, S(O)_{2}R^{12},
 5
            S(0)NR^{12}R^{13}, S(0)_2NR^{12}R^{13}, NR^{14}S(0)R^{12}, NR^{14}S(0)_2R^{12},
           NR^{12}C(0)R^{15}, NR^{12}C(0)OR^{15}, NR^{12}S(0)_2R^{15}, and
           NR12C(O)NHR15:
     R11 is selected from
10
           H, halo, -CF3, -CN, -NO2, C1-6 alkyl,
           C2-6 alkenyl, C2-6 alkynyl, C1-4 haloalkyl, C1-6 alkoxy,
                  C3-10 cycloalkyl,
           C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>,
           arvl substituted with 0-5 R33.
15
            5-10 membered heterocyclic ring system containing from
                  1-4 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
                  R31;
20
           OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13},
                  NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, OC(0)OR^{12}.
                  CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}.
                  S(0)_2R^{12}, S(0)NR^{12}R^{13}, S(0)_2NR^{12}R^{13}, NR^{14}S(0)R^{12},
                  and NR14S(O)2R12;
25
     \mathbb{R}^{12}, at each occurrence, is independently selected from
           C<sub>1-4</sub> alkyl substituted with 0-1 R<sup>12a</sup>,
            C2-4 alkenyl substituted with 0-1 R12a,
           C2-4 alkynyl substituted with 0-1 R12a,
30
           C3-6 cycloalkyl substituted with 0-3 R33,
           aryl substituted with 0-5 R33;
           C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>, and
            5-10 membered heterocyclic ring system containing from
                  1-4 heteroatoms selected from the group
35
                  consisting of N, O, and S substituted with 0-3
                  R31.
```

- R^{12a} , at each occurrence, is independently selected from phenyl substituted with 0-5 R^{33} ; C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R^{31} ;
- R¹³, at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
 - alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or -N(R^{14})-;
- alternatively, R¹² and R¹³ when attached to N may be combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group consisting of N, O, and S, wherein said bicyclic heterocyclic ring system is unsaturated or partially saturated, wherein said bicyclic heterocyclic ring system is substituted with 0-3 R¹⁶;
- R¹⁴, at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;
 - R^{15} , at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
- 30 R¹⁶, at each occurrence, is independently selected from H, OH, F, Cl, CN, NO₂, CF₃, SO₂R⁴⁵, NR⁴⁶R⁴⁷, -C(=O)H, methyl, ethyl, methoxy, ethoxy, trifluoromethyl, trifluoromethoxy, and =O;
- 35 R^{31} , at each occurrence, is independently selected from H, OH, halo, CF₃, SO_2R^{45} , $NR^{46}R^{47}$, C_{1-4} alkyl, and =0;

```
\mathbb{R}^{33}, at each occurrence, is independently selected from
              H, OH, halo, CN, NO<sub>2</sub>, CF<sub>3</sub>, SO<sub>2</sub>R<sup>45</sup>, NR<sup>46</sup>R<sup>47</sup>, -C(=O)H,
              =0, phenyl, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl,
              C3-6 cycloalkyl, C1-4 haloalkyl, C1-4 haloalkyl-oxy-,
             C_{1-4} alkyloxy-, C_{1-4} alkylthio-, C_{1-4} alkyl-C(=0)-,
             C_{1-4} alkyl-C(=0)NH-, C_{1-4} alkyl-OC(=0)-,
             C_{1-4} alkyl-C(=0)0-, C_{3-6} cycloalkyl-oxy-.
             C<sub>3-6</sub> cycloalkylmethyl-oxy-;
             C1-6 alkyl substituted with OH, methoxy, ethoxy,
                propoxy, butoxy, -SO<sub>2</sub>R<sup>45</sup>, -NR<sup>46</sup>R<sup>47</sup>, NR<sup>46</sup>R<sup>47</sup>C(=0)-, or
10
                 (C_{1-4} \text{ alkyl})CO_{2-}; and
             C2-6 alkenyl substituted with OH, methoxy, ethoxy,
                propoxy, butoxy, -SO<sub>2</sub>R<sup>45</sup>, -NR<sup>46</sup>R<sup>47</sup>, NR<sup>46</sup>R<sup>47</sup>C(=O)-, or
                (C_{1-4} \text{ alkyl}) CO_2-;
15
      R^{41}, at each occurrence, is independently selected from
             H, CF<sub>3</sub>, halo, OH, CO<sub>2</sub>H, SO<sub>2</sub>R<sup>45</sup>, NR<sup>46</sup>R<sup>47</sup>, NO<sub>2</sub>, CN,
             C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl
             C_{1-4} alkyl substituted with 0-1 R^{43},
20
             aryl substituted with 0-3 R42, and
             5-10 membered heterocyclic ring system containing from
                    1-4 heteroatoms selected from the group
                    consisting of N, O, and S substituted with 0-3
                    R44;
25
      \mathbb{R}^{42}, at each occurrence, is independently selected from
             H, CF_3, halo, OH, CO_2H, SO_2R^{45}, NR^{46}R^{47}, NO_2, CN,
                   CH (=NH) NH2, NHC (=NH) NH2,
            C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl,
30
                   C3-6 cycloalkyl.
             C<sub>1-4</sub> alkyl substituted with 0-1 R<sup>43</sup>,
             aryl substituted with 0-3 R44, and
            5-10 membered heterocyclic ring system containing from
                   1-4 heteroatoms selected from the group
35
                   consisting of N, O, and S substituted with 0-3
```

R44:

```
R^{43} is C_{3-6} cycloalkyl or aryl substituted with 0-3 R^{44};
      {\bf R}^{44}, at each occurrence, is independently selected from H,
             halo, -OH, NR^{46}R^{47}, CO_2H, SO_2R^{45}, -CF<sub>3</sub>, -OCF<sub>3</sub>, -CN, -
  5
             NO_2, C_{1-4} alkyl, and C_{1-4} alkoxy;
      R^{45} is C_{1-4} alkyl;
      \ensuremath{\mathrm{R}}^{46}, at each occurrence, is independently selected from \ensuremath{\mathrm{H}}
10
             and C_{1-4} alkyl;
      \ensuremath{\text{R}^{47}}\xspace , at each occurrence, is independently selected from H
            and C_{1-4} alkyl;
    n is 1 or 2;
      m is 1 or 2; and
      n plus m is 2, 3, or 4:
      provided when n is 1, m is 2, and \mathbb{R}^7, \mathbb{R}^8, and \mathbb{R}^9 are
      independently selected from H, halogen, C_{1-4} alkyl, C_{1-4}
      alkoxy, C_{1-4} alkylthio or trifluoromethyl; then X is not a
      bond.
      4. A compound of Claim 2 wherein:
25
      X is a bond, -CH_2-, -O-, -S-, -OCH_2-, or -SCH_2-;
      R1 is selected from
             Η.
30
             C_{1-4} alkyl,
             C2-4 alkenyl,
             C_{2-4} alkynyl,
             C3-4 cycloalkyl,
             C_{1-3} alkyl substituted with 0-1 R^2,
35
             C_{2-3} alkenyl substituted with 0-1 \mathbb{R}^2, and
             C_{2-3} alkynyl substituted with 0-1 R^2:
```

```
\mathbb{R}^2, at each occurrence, is independently selected from
            C_{1-4} alkyl,
            C2-4 alkenyl,
            C2-4 alkynyl,
            C3-6 cycloalkyl,
            phenyl substituted with 0-5 R42:
            \text{C}_{3\text{-}6} carbocyclic residue substituted with 0-3 \text{R}^{41}, and
            5-6 membered heterocyclic ring system containing 1, 2,
                  or 3 heteroatoms selected from the group
10
                  consisting of N, O, and S substituted with 0-3
                  R41:
      R<sup>6a</sup> is H, methyl, ethyl, propyl, or butyl;
15
      R6b is H:
      alternatively, R^{6a} and R^{6b} are taken together to form =0 or
            =S:
20
      \ensuremath{\mbox{R}^{7}} and \ensuremath{\mbox{R}^{9}}, at each occurrence, are independently selected
            from
            H, halo, -CF3, -OCF3, -OH, -CN, -NO2, -NR46R47,
            C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,
                  C_{1-4} alkoxy, (C_{1-4} haloalkyl)oxy,
25
            C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>,
            C<sub>1-4</sub> alkyl substituted with 0-2 R<sup>11</sup>.
            C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
            aryl substituted with 0-5 R33, and
           5-6 membered heterocyclic ring system containing 1, 2,
30
                  or 3 heteroatoms selected from the group
                 consisting of N, O, and S substituted with 0-3
                 R31;
     R8 is selected from
35
           H, halo, -CF_3, -OCF_3, -OH, -CN, -NO_2,
           C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,
```

C1-4 alkoxy, (C1-4 haloalkyl)oxy,

```
C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>,
            C<sub>1-4</sub> alkyl substituted with 0-2 R<sup>11</sup>.
            C_{2-4} alkenyl substituted with 0-2 R^{11},
            C2-4 alkynyl substituted with 0-1 R11,
  5
            C<sub>3-10</sub> carbocyclic residue substituted with 0-3 R<sup>33</sup>
            aryl substituted with 0-5 R33.
            5-6 membered heterocyclic ring system containing 1, 2,
                  or 3 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
10
                  R31.
            OR12, SR12, NR12R13, NR12C(O)R15, NR12C(O)OR15.
                  NR^{12}S(0)_2R^{15}, NR^{12}C(0)NHR^{15}, NR^{14}C(0)R^{12}.
                  NR^{14}C(0)OR^{12}, and NR^{14}S(0)_2R^{12};
15
     R11 is selected from
            H, halo, -CF3, -CN, -NO2,
            C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,
                  C_{1-4} alkoxy, (C_{1-4} haloalkyl)oxy,
           C_{3-10} cycloalkyl substituted with 0-2 R^{33},
20
           C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
           aryl substituted with 0-5 R33, and
            5-6 membered heterocyclic ring system containing 1, 2,
                  or 3 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
25
                 p31,
     \mathbb{R}^{12}, at each occurrence, is independently selected from
           C<sub>1-4</sub> alkyl substituted with 0-1 R<sup>12a</sup>,
           C_{2-4} alkenyl substituted with 0-1 R^{12a},
30
           C2-4 alkynyl substituted with 0-1 R12a.
           C<sub>3-6</sub> cycloalkyl substituted with 0-3 R<sup>33</sup>,
           aryl substituted with 0-5 R33;
           C_{3-10} carbocyclic residue substituted with 0-3 R^{33}, and
           5-10 membered heterocyclic ring system containing from
35
                 1-4 heteroatoms selected from the group
                 consisting of N, O, and S substituted with 0-3 \,
                 R31:
```

35

- R^{12a} , at each occurrence, is independently selected from phenyl substituted with 0-5 R^{33} ; C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R^{31} ;
- 10 R¹³, at each occurrence, is independently selected from H, C₁₋₄ alkyl, C₂₋₄ alkenyl, and C₂₋₄ alkynyl;
 - alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or -N(R^{14})-;
 - alternatively, R¹² and R¹³ when attached to N may be combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group consisting of one N, two N, three N, one N one O, and one N one S; wherein said bicyclic heterocyclic ring system is unsaturated or partially saturated, wherein said bicyclic heterocyclic ring system is substituted with 0-2 R¹⁶;
 - R¹⁴, at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;
- R¹⁵, at each occurrence, is independently selected from H, 30 methyl, ethyl, propyl, and butyl;
 - R^{16} , at each occurrence, is independently selected from H, OH, F, Cl, CN, NO₂, methyl, ethyl, methoxy, trifluoromethyl, and trifluoromethoxy;
 - R³¹, at each occurrence, is independently selected from H, OH, halo, CF₃, methyl, ethyl, and propyl;

35

1. 14

- 10 C_{1-6} alkyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, $-SO_2R^{45}$, $-NR^{46}R^{47}$, $NR^{46}R^{47}C$ (=O)-, or $(C_{1-4}$ alkyl) CO_2 -; and
 - C_{2-6} alkenyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, -SO_2R^{45}, -NR^{46}R^{47}, NR^{46}R^{47}C (=0) -, or $(C_{1-4} \text{ alkyl}) \text{CO}_2\text{-};$
 - $\rm R^{41},$ at each occurrence, is independently selected from H, CF3, halo, OH, CO₂H, SO₂R⁴⁵, NR⁴⁶R⁴⁷, NO₂, CN, C₂₋₄ alkenyl, C₂₋₄ alkynyl, C₁₋₃ alkoxy, C₁₋₃ haloalkyl, and C₁₋₃ alkyl;
 - R^{42} , at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO_2R^{45} , $NR^{46}R^{47}$, NO_2 , CN, CH (=NH) NH_2 , NHC (=NH) NH_2 ,
- - R^{43} is cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, phenyl, or pyridyl, each substituted with 0-3 R^{44} ;
 - R⁴⁴, at each occurrence, is independently selected from H, halo, -OH, NR⁴⁶R⁴⁷, CO₂H, SO₂R⁴⁵, -CF₃, -OCF₃, -CN, -NO₂, methyl, ethyl, propyl, butyl, methoxy, ethoxy, propoxy, and butoxy;
 - R45 is methyl, ethyl, propyl, or butyl;

```
15
```

```
R^{46}, at each occurrence, is independently selected from H,
           methyl, ethyl, propyl, and butyl;
     R^{47}, at each occurrence, is independently selected from
 5
           from H, methyl, ethyl, propyl, and butyl;
     n is 1 or 2;
     m is 1 or 2; and
     n plus m is 2 or 3:
10
     provided when n is 1, m is 2, and R^7, R^8, and R^9 are
     independently selected from H, halogen, C_{1-4} alkyl, C_{1-4}
     alkoxy, C1-4 alkylthic or trifluoromethyl; then X is not a
     bond.
15
     5. A compound of Claim 2 wherein:
     X is a bond, -CH2-, -O-, -S-, -OCH2-, or -SCH2-;
     R1 is selected from
            Η,
            C_{1-4} alkyl,
            C2-4 alkenyl,
25
            C2-4 alkynyl,
           C3-4 cycloalkyl,
            C_{1-3} alkyl substituted with 0-1 R^2,
           C_{2-3} alkenyl substituted with 0-1 R^2, and
            C_{2-3} alkynyl substituted with 0-1 R^2;
30
     \mathbb{R}^2, at each occurrence, is independently selected from
          C_{1-4} alkyl,
          C2-4 alkenyl,
          C_{2-4} alkynyl,
35
          C3-6 cycloalkyl.
          phenyl substituted with 0-5 R42;
          C_{3-6} carbocyclic residue substituted with 0-3 \mathbb{R}^{41}, and
```

```
5-6 membered heterocyclic ring system containing 1, 2,
                  or 3 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
                  R41;
  5
      R6a is H;
      R6b is H;
10
      alternatively, R6a and R6b are taken together to form =0;
      {\ensuremath{\mathtt{R}}}^7 and {\ensuremath{\mathtt{R}}}^9, at each occurrence, are independently selected
            from
            H, F, Cl, -CH3, -OCH3, -CF3, -OCF3, -CN, and -NO2,
15
      R8 is selected from
            H, F, Cl, Br, -CF3, -OCF3, -OH, -CN, -NO2,
            C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,
                  C_{1-4} alkoxy, (C_{1-4} haloalkyl)oxy,
20
            C<sub>3-10</sub> cycloalkyl substituted with 0-2 R<sup>33</sup>.
            C_{1-4} alkyl substituted with 0-2 R^{11},
            C2-4 alkenyl substituted with 0-2 R11,
            C2-4 alkynyl substituted with 0-1 R11.
            C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
25
            aryl substituted with 0-5 R33.
            5-6 membered heterocyclic ring system containing 1, 2,
                  or 3 heteroatoms selected from the group
                  consisting of N, O, and S substituted with 0-3
                  R31.
           OR^{12}, SR^{12}, NR^{12}R^{13}, NR^{12}C(O)R^{15}, NR^{12}C(O)OR^{15},
30
                  NR^{12}S(0)_2R^{15}, NR^{12}C(0)NHR^{15}, NR^{14}C(0)R^{12}.
                  NR^{14}C(0)OR^{12}, and NR^{14}S(0) \circ R^{12}.
     R11 is selected from
35
           H, halo, -CF3, -CN, -NO2,
           C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,
                 C_{1-4} alkoxy, (C_{1-4} haloalkyl)oxy,
```

25

5

10

C₃₋₁₀ cycloalkyl substituted with 0-2 R³³. C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , aryl substituted with 0-5 R33, and 5-6 membered heterocyclic ring system containing 1, 2, or 3 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3R31.

 \mathbb{R}^{12} , at each occurrence, is independently selected from C₁₋₄ alkyl substituted with 0-1 R^{12a}, C2-4 alkenyl substituted with 0-1 R12a. C2-4 alkynyl substituted with 0-1 R12a. C₃₋₆ cycloalkyl substituted with 0-3 R³³, aryl substituted with 0-5 R33; 15 C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 $\,$ R31;

 R^{12a} , at each occurrence, is independently selected from phenyl substituted with 0-5 R33; C_{3-10} carbocyclic residue substituted with 0-3 \mathbb{R}^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 $\,$ R31:

 \mathbb{R}^{13} , at each occurrence, is independently selected from 30 H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;

alternatively, \mathbf{R}^{12} and \mathbf{R}^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or $-N(R^{14})$ -;

alternatively, ${\bf R}^{12}$ and ${\bf R}^{13}$ when attached to N may be 35 combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3

20

25

35

5

heteroatoms selected from the group consisting of N, O, and S; wherein said bicyclic heterocyclic ring system is selected from indolyl, indolinyl, indazolyl, benzimidazolyl, benzimidazolyl, benzimidazolyl, benzimidazolyl, duinolinyl, tetrahydroquinolinyl, isoquinolinyl, and tetrahydroisoquinolinyl; wherein said bicyclic heterocyclic ring system is substituted with 0-1 \mathbb{R}^{16} ;

- R¹⁴, at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;
 - R¹⁵, at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;
 - R¹⁶, at each occurrence, is independently selected from H, OH, F, Cl, CN, NO₂, methyl, ethyl, methoxy, ethoxy, trifluoromethyl, and trifluoromethoxy;
 - $R^{31},\ \mbox{at each occurrence, is independently selected from H, OH, halo, CF_3, methyl, ethyl, and propyl;} \label{eq:R31}$
 - R³³, at each occurrence, is independently selected from H, OH, halo, CN, NO₂, CF₃, SO₂R⁴⁵, NR⁴⁶R⁴⁷, -C(=O)H, phenyl, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-6} cycloalkyl, C_{1-4} haloalkyl, C_{1-4} haloalkyl-oxy-, C_{1-4} alkyloxy-, C_{1-4} alkylthio-, C_{1-4} alkyl-C(=O) NH-, C_{1-4} alkyl-OC(=O)-, C_{1-4} alkyl-C(=O)O-, C_{3-6} cycloalkyl-oxy-,
- C₃₋₆ cycloalkylmethyl-oxy-;

 C₁₋₆ alkyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, -So₂R⁴⁵, -NR⁴⁶R⁴⁷, NR⁴⁶R⁴⁷C(=O)-, or (C₁₋₄ alkyl)CO₂-; and
 - $\rm C_{2-6}$ alkenyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, -SO₂R⁴⁵, -NR⁴⁶R⁴⁷, NR⁴⁶R⁴⁷C(=O)-, or (C₁₋₄ alkyl)CO₂-;
 - \mathbb{R}^{41} , at each occurrence, is independently selected from

```
H, CF_3, halo, OH, CO_2H, SO_2R^{45}, NR^{46}R^{47}, NO_2, CN, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-3} alkoxy, C_{1-3} haloalkyl, and C_{1-3} alkyl;
```

5 R⁴², at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO_2R^{45} , $NR^{46}R^{47}$, NO_2 , CN, CH(=NH)NH₂, NHC(=NH)NH₂,

 C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-3} alkoxy, C_{1-3} haloalkyl, C_{3-6} cycloalkyl, and C_{1-3} alkyl;

10

15

 R^{43} is cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, phenyl, or pyridyl, each substituted with 0-3 R^{44} ;

 $\rm R^{44}$, at each occurrence, is independently selected from H, halo, -OH, NR^46R^47, CO_2H, SO_2R^45, -CF_3, -OCF_3, -CN, -NO_2, methyl, ethyl, propyl, butyl, methoxy, ethoxy, propoxy, and butoxy;

R45 is methyl, ethyl, propyl, or butyl;

20

35

 R^{46} , at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;

 R^{47} , at each occurrence, is independently selected from 25 from H, methyl, ethyl, propyl, and butyl;

n is 1; and m is 1.

30 6. A compound of Claim 2 wherein:

X is a bond, $-CH_2-$, -O-, -S-, $-OCH_2-$, or $-SCH_2-$;

R1 is selected from H,

 C_{1-5} alkyl substituted with 0-1 R^2 , C_{2-5} alkenyl substituted with 0-1 R^2 , and C_{2-3} alkynyl substituted with 0-1 R^2 :

```
\mathbb{R}^2 is cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, or
                phenyl;
        R6a is H.
         R6b is H:
         \ensuremath{\mbox{R}^{7}} and \ensuremath{\mbox{R}^{9}}, at each occurrence, are independently selected
    10
                from H, F, Cl, -CH3, -OCH3, -CF3, -OCF3, -CN, and -NO2;
         R^8 is selected from R^{11};
               methyl substituted with R11;
               phenyl substituted with 0-3 R33:
   15
               pyridyl substituted with 0-2 R33:
               OR^{12}, SR^{12}, NR^{12}R^{13}, NR^{12}C(0)R^{15}, NR^{12}C(0)OR^{15},
                     NR^{12}S(0)_2R^{15}, NR^{12}C(0)NHR^{15}, NR^{14}C(0)R^{12}.
                     NR^{14}C(0)OR^{12}, and NR^{14}S(0)_2R^{12};
20
         R11 is selected from
               phenyl- substituted with 0-5 fluoro;
               pyridyl substituted with 0-2 R33:
               naphthyl- substituted with 0-2 R33;
               2-(H3CCH2C(=0))-phenyl- substituted with R33;
   25
               2-(H<sub>3</sub>CC(=0))-phenyl- substituted with R<sup>33</sup>;
               2-(HC(=0))-phenyl- substituted with R33.
               2-(H3CCH(OH))-phenyl- substituted with R33:
               2-(H<sub>3</sub>CCH<sub>2</sub>CH(OH))-phenyl- substituted with R<sup>33</sup>;
               2-(HOCH2)-phenyl- substituted with R33;
   30
               2-(HOCH<sub>2</sub>CH<sub>2</sub>)-phenyl- substituted with R<sup>33</sup>;
               2-(H3COCH2)-phenyl- substituted with R33;
               2-(H3COCH2CH2)-phenyl- substituted with R33;
               2-(H3CCH(OMe))-phenyl- substituted with R33;
               2-(H<sub>3</sub>COC(=0))-phenyl- substituted with R<sup>33</sup>;
   35
               2-(HOCH2CH=CH)-phenyl- substituted with R33;
               2-((MeOC=O)CH=CH)-phenyl- substituted with R<sup>33</sup>;
               2-(methyl)-phenyl- substituted with R33;
```

```
2-(ethyl)-phenyl- substituted with R33;
               2-(i-propyl)-phenyl- substituted with R33;
               2-(F3C)-phenyl- substituted with R33:
               2-(NC)-phenyl- substituted with R33.
     5
               2-(H3CO)-phenyl- substituted with R33;
               2-(fluoro)-phenyl- substituted with R33;
               2-(chloro)-phenyl- substituted with R33:
               3-(NC)-phenyl- substituted with R33:
               3-(H3CO)-phenyl- substituted with R33;
              3-(fluoro)-phenyl- substituted with R33:
    10
               3-(chloro)-phenyl- substituted with R33:
               3-(H3C)-phenyl- substituted with R33;
(3
              3-(F<sub>3</sub>C)-phenyl- substituted with R<sup>33</sup>;
State State State
               3-(H2CS)-phenyl- substituted with R33;
    15
              4-(NC)-phenyl- substituted with R33;
100
              4-(fluoro)-phenyl- substituted with R33;
1.1
              4-(chloro)-phenyl- substituted with R33;
10
              4-(H3CS)-phenyl- substituted with R33;
Dane office
              4-(H<sub>3</sub>CO)-phenyl- substituted with R<sup>33</sup>;
   20
13
              4-(ethoxy)-phenyl- substituted with R33:
0
              4-(i-propoxy)-phenyl- substituted with R33:
1
              4-(i-butoxy)-phenyl- substituted with R33;
              4-(H<sub>3</sub>CCH<sub>2</sub>CH<sub>2</sub>C(=O))-phenyl- substituted with R<sup>33</sup>;
              4-((H_3C)<sub>2</sub>CHC(=0))-phenyl- substituted with R^{33};
   25
              4-(H3CCH2C(=0))-phenyl- substituted with R33;
              4-(H<sub>3</sub>CC(=0))-phenyl- substituted with R<sup>33</sup>;
              4-(H3CCH2CH2CH(OH))-phenyl- substituted with R33:
              4-((H<sub>3</sub>C)<sub>2</sub>CHCH(OH))-phenyl- substituted with R<sup>33</sup>;
              4-(H3CCH2CH(OH))-phenyl- substituted with R33;
   30
              4-(H3CCH(OH))-phenyl- substituted with R33;
              4-(cyclopropyloxy)-phenyl- substituted with R33;
              4-(cyclobutyloxy)-phenyl- substituted with R33; and
              4-(cyclopentyloxy)-phenyl- substituted with R33;
   35
        R12 is selected from
              methyl substituted with R11.
              phenyl substituted with 0-5 fluoro;
```

```
pyridyl substituted with 0-2 R33;
                naphthyl substituted with 0-2 R33:
                2-(H<sub>3</sub>CCH<sub>2</sub>C(=O))-phenyl- substituted with R<sup>33</sup>;
                2-(H<sub>3</sub>CC(=0))-phenyl- substituted with R<sup>33</sup>;
     5
                2-(HC(=O))-phenyl- substituted with R33:
                2-(H<sub>3</sub>CCH(OH))-phenyl- substituted with R<sup>33</sup>;
                2-(H<sub>3</sub>CCH<sub>2</sub>CH(OH))-phenyl- substituted with R<sup>33</sup>;
                2-(HOCH<sub>2</sub>)-phenyl- substituted with R<sup>33</sup>;
                2-(HOCH<sub>2</sub>CH<sub>2</sub>)-phenyl- substituted with R<sup>33</sup>:
    10
                2-(H3COCH2)-phenyl- substituted with R33;
               2-(H3COCH2CH2)-phenyl- substituted with R33;
               2-(H<sub>3</sub>CCH(OMe))-phenyl- substituted with R<sup>33</sup>:
               2-(H<sub>3</sub>COC(=O))-phenyl- substituted with R<sup>33</sup>;
               2-(HOCH2CH=CH)-phenyl- substituted with R33;
200
               2-((MeOC=O)CH=CH)-phenyl- substituted with R<sup>33</sup>;
    15
               2-(methyl)-phenyl- substituted with R33;
13
               2-(ethyl)-phenyl- substituted with R33;
Li
               2-(i-propyl)-phenyl- substituted with R33:
               2-(F<sub>3</sub>C)-phenyl- substituted with R<sup>33</sup>;
70
100
    20
               2-(NC)-phenyl- substituted with R33;
100
               2-(H3CO)-phenyl- substituted with R33;
1.4
               2-(fluoro)-phenyl- substituted with R33;
               2-(chloro)-phenyl- substituted with R33;
               3-(NC)-phenyl- substituted with R33;
    25
               3-(H3CO)-phenyl- substituted with R33;
               3-(fluoro)-phenyl- substituted with R33;
               3-(chloro)-phenyl- substituted with R33;
               3-(H_3C)-phenyl- substituted with R^{33};
               3-(F3C)-phenyl- substituted with R33:
   30
               3-(H<sub>3</sub>CS)-phenyl- substituted with R<sup>33</sup>;
               4-(fluoro)-phenyl- substituted with R33;
               4-(chloro)-phenyl- substituted with R33;
              4-(H<sub>3</sub>CS)-phenyl- substituted with R<sup>33</sup>;
               4-(H<sub>3</sub>CO)-phenyl- substituted with R<sup>33</sup>;
   35
              4-(ethoxy)-phenyl- substituted with R33;
              4-(i-propoxy)-phenyl- substituted with R33:
              4-(i-butoxy)-phenyl- substituted with R33;
```

1

29

į, į

10

20

30

4-(H3CCH2CH2C(=0))-phenyl- substituted with R33: 4-((H₃C)₂CHC(=0))-phenyl- substituted with R³³: 4-(H3CCH2C(=0))-phenyl- substituted with R33; 4-(H3CC(=0))-phenyl- substituted with R33; 4-(H₃CCH₂CH₂CH(OH))-phenyl- substituted with R³³; 4-((H₃C)₂CHCH(OH))-phenyl- substituted with R³³; 4-(H₃CCH₂CH(OH))-phenyl- substituted with R³³; 4-(H3CCH(OH))-phenyl- substituted with R33; 4-(cyclopropyloxy)-phenyl- substituted with R33; 4-(cyclobutyloxy)-phenyl- substituted with R33; and

4-(cyclopentyloxy)-phenyl- substituted with R33;

- R13 is H, methyl, or ethyl;
- alternatively, \mathbb{R}^{12} and \mathbb{R}^{13} join to form a 5- or 6-membered 15 ring selected from pyrrolyl, pyrrolidinyl, imidazolyl, piperidinyl, piperizinyl, methylpiperizinyl, and morpholinyl;
- alternatively, ${\bf R}^{12}$ and ${\bf R}^{13}$ when attached to N may be combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group consisting of N, O, and S; wherein said bicyclic heterocyclic ring system is selected from indolyl, indolinyl, indazolyl, 25 benzimidazolyl, benzimidazolinyl, benztriazolyl, quinolinyl, tetrahydroquinolinyl, isoquinolinyl, and tetrahydroisoquinolinyl; wherein said bicyclic heterocyclic ring system is substituted with 0-1 R^{16} ;
 - R15 is H, methyl, ethyl, propyl, or butyl;
- \mathbb{R}^{16} , at each occurrence, is independently selected from ${\tt H}, {\tt OH}, {\tt F}, {\tt Cl}, {\tt CN}, {\tt NO}_2, {\tt methyl}, {\tt ethyl}, {\tt methoxy}, {\tt ethoxy},$ 35 trifluoromethyl, and trifluoromethoxy;
 - \mathbb{R}^{33} , at each occurrence, is independently selected from

15

30

H, F, Cl, -CH₃, -OCH₃, -SCH₃, -CF₃, -OCF₃, -CN, and -NO₂:

n is 1; and m is 1.

A compound of Claim 2 of Formula (I-a)

wherein:

b is a single bond wherein the bridging hydrogens are either cis or trans;

X is a bond, $-CH_2-$, -O-, -S-, $-OCH_2-$, or $-SCH_2-$;

R1 is selected from

20 hydrogen, methyl, ethyl, n-propyl, n-butyl, s-butyl, t-butyl, n-pentyl, n-hexyl, 2-propyl, 2-butyl, 2-pentyl, 2-hexyl, 2-methylpropyl, 2-methylbutyl, 2-methylpentyl, 2-ethylbutyl, 3-methylpentyl, 3-methylbutyl, 4-methylpentyl, 2-fluoroethyl, 2,2-difluoroethyl, 2,2,2-trifluoroethyl,

2-propenyl, 2-methyl-2-propenyl, trans-2-butenyl, 3-methyl-2-butenyl, 3-butenyl, trans-2-pentenyl, cis-2-pentenyl, 4-pentenyl, 4-methyl-3-pentenyl, 3,3-dichloro-2-propenyl, trans-3-phenyl-2-propenyl,

cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, cyclopropylmethyl, cyclobutylmethyl, cyclopentylmethyl, cyclohexylmethyl,

```
1.14
10
(1)
[]
[] 15
17
17
£
14
```

```
benzyl, 2-methylbenzyl, 3-methylbenzyl, 4-methylbenzyl,
         2,5-dimethylbenzyl, 2,4-dimethylbenzyl, 3,5-
         dimethylbenzyl,
        2,4,6-trimethyl-benzyl, 3-methoxy-benzyl, 3,5-dimethoxy-
  5
        benzyl, pentafluorobenzyl, 2-phenylethyl, 1-phenyl-2-
        propyl, 4-phenylbutyl, 4-phenylbenzyl, 2-phenylbenzyl,
         (2,3-dimethoxy-phenyl)C(=0)-, (2,5-dimethoxy-
10
        phenyl)C(=0)-, (3,4-dimethoxy-phenyl)C(=0)-,
        (3,5-dimethoxy-phenyl)C(=0)-, cyclopropyl-C(=0)-,
        isopropyl-C(=0)-, ethyl-CO<sub>2</sub>-, propyl-CO<sub>2</sub>-, t-butyl-CO<sub>2</sub>-,
        2,6-dimethoxy-benzyl, 2,4-dimethoxy-benzyl,
        2,4,6-trimethoxy-benzyl, 2,3-dimethoxy-benzyl,
        2,4,5-trimethoxy-benzyl, 2,3,4-trimethoxy-benzyl,
        3,4-dimethoxy-benzyl, 3,4,5-trimethoxy-benzyl,
        (4-fluoro-phenyl)ethyl,
        -CH=CH<sub>2</sub>, -CH<sub>2</sub>-CH=CH<sub>2</sub>, -CH=CH-CH<sub>3</sub>, -C=CH, -C=C-CH<sub>3</sub>, and
20
        -CH2-C≡CH; and
     R6a is H:
     R6b is H;
25
     alternatively, R^{6a} and R^{6b} are taken together to form =0:
     \mathbb{R}^7, \mathbb{R}^8, and \mathbb{R}^9, at each occurrence, are independently
           selected from
       hydrogen, fluoro, chloro, bromo, cyano, methyl, ethyl,
30
       propyl, isopropyl, butyl, t-butyl, nitro,
       trifluoromethyl, methoxy, ethoxy, isopropoxy,
       trifluoromethoxy, phenyl;
35
       2-Cl-phenyl; 2-F-phenyl; 2-Br-phenyl; 2-CN-phenyl;
```

```
2-Me-phenyl; 2-CF3-phenyl; 2-MeO-phenyl; 2-CF3O-phenyl;
          2-NO2-phenyl; 2-MeS-phenyl; 2-CHO-phenyl; 2-HOCH2-
          phenyl;
    5
          3-Cl-phenyl; 3-F-phenyl; 3-Br-phenyl; 3-CN-phenyl;
          3-Me-phenyl; 3-Et-phenyl; 3-n-Pr-phenyl; 3-isoPr-phenyl;
          3-n-Bu-phenyl; 3-CF3-phenyl; 3-MeO-phenyl; 3-MeS-phenyl;
          3-isopropoxyphenyl; 3-CF3O-phenyl; 3-NO2-phenyl;
          3-CHO-phenyl; 3-HOCH2-phenyl; 3-MeOCH2-phenyl;
   10
          3-Me2NCH2-phenyl;
          4-Cl-phenyl; 4-F-phenyl; 4-Br-phenyl; 4-CN-phenyl;
          4-Me-phenyl; 4-Et-phenyl; 4-n-Pr-phenyl;
15
          4-iso-Pr-phenyl; 4-n-Bu-phenyl; 4-CF3-phenyl;
          4-MeO-phenyl; 4-isopropoxyphenyl; 4-CF3O-phenyl;
          4-MeS-phenyl;
          4-acetylphenyl; 3-acetamidophenyl; 4-pyridyl;
          2-furanyl; 2-thiophenyl; 2-naphthyl; 1-pyrrolidinyl,
   20
          2,3-diCl-phenyl; 2,3-diF-phenyl; 2,3-diMe-phenyl;
          2,3-diCF3-phenyl; 2,3-diMeO-phenyl; 2,3-diCF3O-phenyl;
          2,4-diCl-phenyl; 2,4-diF-phenyl; 2,4-diMe-phenyl;
   25
          2,4-diCF3-phenyl; 2,4-diMeO-phenyl; 2,4-diCF3O-phenyl;
          2,5-diCl-phenyl; 2,5-diF-phenyl; 2,5-diMe-phenyl;
          2,5-diCF3-phenyl; 2,5-diMeO-phenyl; 2,5-diCF3O-phenyl;
   30
          2,6-diCl-phenyl; 2,6-diF-phenyl; 2,6-diMe-phenyl;
          2,6-diCF3-phenyl; 2,6-diMeO-phenyl; 2,6-diCF3O-phenyl;
          3,4-diCl-phenyl; 3,4-diF-phenyl; 3,4-diMe-phenyl;
         3,4-diCF3-phenyl; 3,4-diMeO-phenyl; 3,4-diCF3O-phenyl;
  35
         2,4,6-triCl-phenyl; 2,4,6-triF-phenyl;
         2,4,6-triMe-phenyl; 2,4,6-triCF3-phenyl;
```

```
and the first term and after print term to the same and the same term to the same term of the term terms.
```

```
2,4,6-triMeO-phenyl; 2,4,6-triCF3O-phenyl;
        2,4,5-triMe-phenyl; 2,3,4-triF-phenyl;
        2-Me-4-MeO-5-F-phenyl; 2,6-diCl-4-MeO-phenyl;
        2,4-diMeO-6-F-phenyl; 2,6-diF-4-Cl-phenyl;
        2,3,4,6-tetraF-phenyl; 2,3,4,5,6-pentaF-phenyl;
  5
        2-Cl-4-F-phenyl; 2-Cl-6-F-phenyl; 2-Cl-3-Me-phenyl;
        2-Cl-4-MeO-phenyl; 2-Cl-4-EtO-phenyl;
        2-Cl-4-iPrO-phenyl; 2-Cl-4-CF3-phenyl;
 10
        2-C1-4-CF3O-phenyl; 2-C1-4-(CHF2)O-phenyl;
        2-F-3-Cl-phenyl; 2-F-4-MeO-phenyl; 2-F-5-Me-phenyl;
        2-Me-3-Cl-phenyl; 2-Me-3-CN-phenyl; 2-Me-4-Cl-phenyl;
        2-Me-4-F-phenyl; 2-Me-4-CN-phenyl; 2-Me-4-MeO-phenyl;
15
        2-Me-4-EtO-phenyl; 2-Me-4-MeS-phenyl;
        2-Me-4-H2NCO-phenyl; 2-Me-4-MeOC(=0)-phenyl;
        2-Me-4-CH<sub>3</sub>C(=0)-phenyl; 2-Me-5-F-phenyl;
        2-Et-4-MeO-phenyl; 2-MeO-5-F-phenyl:
        2-MeO-4-isopropyl-phenyl; 2-CF3-4-Cl-phenyl;
20
        2-CF3-4-F-phenyl; 2-CF3-4-MeO-phenyl;
        2-CF3-4-EtO-phenyl; 2-CF3-4-iPrO-phenyl;
        2-CF<sub>3</sub>-4-CN-phenyl; 2-CF<sub>3</sub>-6-F-phenyl;
        2-CHO-4-MeO-phenyl; 2-MeOC(=0)-3-MeO-phenyl;
        2-CH<sub>3</sub>CH(OH)-4-MeO-phenyl; 2-CH<sub>3</sub>CH(OH)-4-F-phenyl;
25
        2-CH<sub>3</sub>CH(OH)-4-Cl-phenyl; 2-CH<sub>3</sub>CH(OH)-4-Me-phenyl;
        2-CH_3CH(OMe)-4-MeO-pheny1; 2-CH_3C(=O)-4-MeO-pheny1;
        2-CH<sub>3</sub>C(=0)-4-F-phenyl; 2-CH<sub>3</sub>C(=0)-4-Cl-phenyl;
        2-CH_3C(=0)-4-Me-phenyl; 2-H_2C(OH)-4-MeO-phenyl;
        2-H_2C(OMe)-4-MeO-phenyl; 2-H_3CCH_2CH(OH)-4-MeO-phenyl;
30
        2-H_3CCH_2C(=0)-4-MeO-phenyl; 2-CH_3CO_2CH_2CH_2-4-MeO-phenyl;
        (Z)-2-HOCH2CH=CH-4-MeO-phenyl;
        (E) -2-HOCH2CH=CH-4-MeO-phenyl;
        (Z) -2-CH3CO2CH=CH-4-MeO-phenvl;
        (E) -2-CH3CO2CH=CH-4-MeO-phenyl;
35
       2-CH3OCH2CH2-4-MeO-phenyl;
       3-CN-4-F-phenyl; 3-HoNCO-4-F-phenyl:
```

```
The first that the first and the first are and the first a
```

```
(2-Cl-phenyl)-CH=CH-; (3-Cl-phenyl)-CH=CH-;
        (2,6-diF-phenyl)-CH=CH-; phenyl-CH=CH-;
        (2-Me-4-MeO-phenyl)-CH=CH-;
        cyclohexyl; cyclopentyl; cyclohexylmethyl; benzyl;
        2-F-benzyl; 3-F-benzyl; 4-F-benzyl; 3-MeO-benzyl;
        3-OH-benzyl; 2-MeO-benzyl; 2-OH-benzyl;
        tetrahydroquinolin-1-vl;
        tetrahydroindolin-1-yl;
1.0
        tetrahydroisoindolin-1-vl:
        phenyl-S-; phenyl-NH-; pyrid-3-yl-NH-;
        (4-Me-pyrid-3-yl)-NH-; (4-Cl-pyrid-3-yl)-NH-;
        (1-naphthyl)-NH-; (2-naphthyl)-NH-;
15
        (2-Me-naphth-1-yl)-NH-; (4-Me-naphth-1-yl)-NH-;
        (3-quinolinvl)-NH-:
        (2-[1,1'-biphenyl])-NH-; (3-[1,1'-biphenyl])-NH-;
        (4-[1,1'-biphenyl])-NH-; (2-F-phenyl)-NH-:
20
        (2-Cl-phenyl)-NH-; (2-CF3-phenyl)-NH-;
        (2-CH3-phenyl)-NH-; (2-OMe-phenyl)-NH-;
        (2-CN-phenyl)-NH-; (2-OCF3-phenyl)-NH-;
        (2-SMe-phenyl)-NH-; (3-F-phenyl)-NH-;
       (3-Cl-phenyl)-NH-; (3-CF<sub>3</sub>-phenyl)-NH-;
25
        (3-CH3-phenyl)-NH-; (3-OMe-phenyl)-NH-;
       (3-CN-phenyl)-NH-; (3-OCF3-phenyl)-NH-;
       (3-SMe-phenyl)-NH-; (4-F-phenyl)-NH-;
       (4-Cl-phenyl)-NH-; (4-CF3-phenyl)-NH-;
       (4-CH3-phenyl)-NH-; (4-OMe-phenyl)-NH-;
30
       (4-CN-phenyl)-NH-; (4-OCF3-phenyl)-NH-;
       (4-SMe-phenyl)-NH-; (2,3-diCl-phenyl)-NH-;
       (2,4-diCl-phenyl)-NH-; (2,5-diCl-phenyl)-NH-;
       (2,6-diCl-phenyl)-NH-; (3,4-diCl-phenyl)-NH-;
       (3,5-diCl-phenyl)-NH-; (2,3-diF-phenyl)-NH-;
35
       (2,4-diF-phenyl)-NH-; (2,5-diF-phenyl)-NH-;
       (2,6-diF-phenyl)-NH-; (3,4-diF-phenyl)-NH-;
       (3,5-diF-phenyl)-NH-; (2,3-diCH3-phenyl)-NH-;
```

```
(2,4-diCH3-phenyl)-NH-; (2,5-diCH3-phenyl)-NH-;
            (2,6-diCH3-phenyl)-NH-; (3,4-diCH3-phenyl)-NH-;
            (3,5-diCH3-phenyl)-NH-; (2,3-diCF3-phenyl)-NH-;
            (2,4-diCF3-phenyl)-NH-; (2,5-diCF3-phenyl)-NH-;
     5
            (2,6-diCF3-phenyl)-NH-; (3,4-diCF3-phenyl)-NH-;
            (3,5-diCF3-phenyl)-NH-; (2,3-diOMe-phenyl)-NH-;
            (2,4-dioMe-phenyl)-NH-; (2,5-dioMe-phenyl)-NH-;
            (2,6-diOMe-phenyl)-NH-; (3,4-diOMe-phenyl)-NH-;
            (3,5-diOMe-phenyl)-NH-; (2-F-3-Cl-phenyl)-NH-;
    10
            (2-F-4-Cl-phenyl)-NH-; (2-F-5-Cl-phenyl)-NH-;
            (2-F-6-C1-phenyl)-NH-; (2-F-3-CH3-phenyl)-NH-;
            (2-F-4-CH3-phenyl)-NH-; (2-F-5-CH3-phenyl)-NH-;
la di
0
            (2-F-6-CH3-phenyl)-NH-; (2-F-3-CF3-phenyl)-NH-;
11
            (2-F-4-CF3-phenyl)-NH-; (2-F-5-CF3-phenyl)-NH-;
£75
   15
            (2-F-6-CF3-phenyl)-NH-; (2-F-3-OMe-phenyl)-NH-;
1.7
            (2-F-4-OMe-phenyl)-NH-; (2-F-5-OMe-phenyl)-NH-;
13
U
            (2-F-6-OMe-phenyl)-NH-; (2-Cl-3-F-phenyl)-NH-;
            (2-Cl-4-F-phenyl)-NH-; (2-Cl-5-F-phenyl)-NH-;
L
1
            (2-Cl-6-F-phenyl)-NH-; (2-Cl-3-CH3-phenyl)-NH-;
   20
            (2-Cl-4-CH3-phenyl)-NH-; (2-Cl-5-CH3-phenyl)-NH-;
1
           (2-Cl-6-CH3-phenyl)-NH-; (2-Cl-3-CF3-phenyl)-NH-;
100
           (2-Cl-4-CF3-phenyl)-NH-; (2-Cl-5-CF3-phenyl)-NH-;
           (2-Cl-6-CF3-phenyl)-NH-; (2-Cl-3-OMe-phenyl)-NH-;
           (2-Cl-4-OMe-phenyl)-NH-; (2-Cl-5-OMe-phenyl)-NH-;
   25
           (2-Cl-6-OMe-phenyl)-NH-; (2-CH<sub>3</sub>-3-F-phenyl)-NH-;
           (2-CH<sub>3</sub>-4-F-phenyl)-NH-; (2-CH<sub>3</sub>-5-F-phenyl)-NH-;
           (2-CH<sub>3</sub>-6-F-phenyl)-NH-; (2-CH<sub>3</sub>-3-Cl-phenyl)-NH-;
           (2-CH<sub>3</sub>-4-Cl-phenyl)-NH-; (2-CH<sub>3</sub>-5-Cl-phenyl)-NH-;
           (2-CH<sub>3</sub>-6-Cl-phenyl)-NH-; (2-CH<sub>3</sub>-3-CF<sub>3</sub>-phenyl)-NH-;
   30
           (2-CH3-4-CF3-phenyl)-NH-; (2-CH3-5-CF3-phenyl)-NH-;
           (2-CH<sub>3</sub>-6-CF<sub>3</sub>-phenyl)-NH-; (2-CH<sub>3</sub>-3-OMe-phenyl)-NH-;
           (2-CH3-4-OMe-phenyl)-NH-; (2-CH3-5-OMe-phenyl)-NH-;
           (2-CH<sub>3</sub>-6-OMe-phenyl)-NH-; (2-CF<sub>3</sub>-3-F-phenyl)-NH-;
           (2-CF3-4-F-phenyl)-NH-; (2-CF3-5-F-phenyl)-NH-;
   35
           (2-CF3-6-F-phenyl)-NH-; (2-CF3-3-Cl-phenyl)-NH-;
           (2-CF<sub>3</sub>-4-Cl-phenyl)-NH-; (2-CF<sub>3</sub>-5-Cl-phenyl)-NH-;
```

(2-CF₃-6-Cl-phenyl)-NH-; (2-CF₃-3-CH₃-phenyl)-NH-;

```
(2-CF<sub>3</sub>-4-CH<sub>3</sub>-phenyl)-NH-; (2-CH<sub>3</sub>-5-CF<sub>3</sub>-phenyl)-NH-;
            (2-CF<sub>3</sub>-6-CH<sub>3</sub>-phenyl)-NH-; (2-CF<sub>3</sub>-3-OMe-phenyl)-NH-;
            (2-CF<sub>3</sub>-4-OMe-phenyl)-NH-; (2-CF<sub>3</sub>-5-OMe-phenyl)-NH-;
            (2-CF<sub>3</sub>-6-OMe-phenyl)-NH-; (2-OMe-3-F-phenyl)-NH-;
            (2-OMe-4-F-phenyl)-NH-; (2-OMe-5-F-phenyl)-NH-;
     5
            (2-OMe-6-F-phenyl)-NH-; (2-OMe-3-Cl-phenyl)-NH-;
            (2-OMe-4-Cl-phenyl)-NH-; (2-OMe-5-Cl-phenyl)-NH-;
            (2-OMe-6-Cl-phenyl)-NH-; (2-OMe-4-CN-phenyl)-NH-;
            (2-OMe-4-CHO-phenyl)-NH-; (2-OMe-3-CH3-phenyl)-NH-;
    10
            (2-OMe-4-CH3-phenyl)-NH-; (2-OMe-5-CH3-phenyl)-NH-;
            (2-OMe-6-CH3-phenyl)-NH-; (2-OMe-3-CF3-phenyl)-NH-;
            (2-OMe-4-CF3-phenyl)-NH-; (2-OMe-5-CF3-phenyl)-NH-;
            (2-OMe-6-CF3-phenyl)-NH-; (2-acetyl-4-Cl-phenyl)-NH-;
()
100
            (2-acetyl-4-Me-phenyl)-NH-; (2-acetyl-4-MeO-phenyl)-NH-;
   15
            (2-CH3CH(OH)-4-Cl-phenyl)-NH-;
(1)
            (2-CH3CH(OH)-4-Me-phenyl)-NH-;
100
            (2-CH<sub>3</sub>CH (OH) -4-MeO-phenyl) -NH-;
U
Bull to the state of
            (3-CF<sub>3</sub>-4-Cl-phenyl)-NH-; (3-F-4-CHO-phenyl)-NH-;
   20
            (3-CH<sub>3</sub>-4-CN-phenyl)-NH-; (3-CH<sub>3</sub>-4-MeO-phenyl)-NH-;
            (3-CH<sub>3</sub>-4-Cl-phenyl)-NH-; (3-CH<sub>3</sub>-4-F-phenyl)-NH-;
            (3-F-5-CF3-phenv1)-NH-:
            (3-CH_3-4-CO_2Me-phenyl)NH-; (3-CF_3-4-C(0)CH_3-phenyl)NH-;
   25
            (3-CHO-4-OMe-phenyl)-NH-; (4-F-3-CF3-phenyl)-NH-;
            (2,3,5-triCl-phenyl)-NH-; (2,4,5-triF-phenyl)-NH-;
            (2,6-diCl-3-Me-phenyl)-NH-; (3,5-diMe-4-MeO-phenyl)-NH-;
            (2-F-3-C1-6-CF3-phenyl)-NH-;
   30
           benzyl-NH-; (3-quinolinyl)CH2NH-; (2-F-phenyl)CH2NH-;
           (2-Cl-phenyl)CH2NH-; (2-CF3-phenyl)CH2NH-;
           (2-CH3-phenyl) CH2NH-; (2-OMe-phenyl) CH2NH-;
           (2-CN-phenyl) CH2NH-; (2-OCF3-phenyl) CH2NH-;
   35
           (2-SMe-phenyl)CH2NH-; (3-F-phenyl)CH2NH-;
           (3-Cl-phenyl) CH2NH-; (3-CF3-phenyl) CH2NH-;
           (3-CH3-phenyl) CH2NH-; (3-OMe-phenyl) CH2NH-;
```

```
. The state of the
```

```
(3-CN-phenyl)CH2NH-; (3-OCF3-phenyl)CH2NH-;
        (3-SMe-phenyl)CH2NH-; (4-F-phenyl)CH2NH-;
        (4-Cl-phenyl) CH2NH-; (4-CF3-phenyl) CH2NH-;
        (4-CH3-phenyl) CH2NH-; (4-OMe-phenyl) CH2NH-;
 5
        (4-CN-phenyl) CH2NH-; (4-OCF3-phenyl) CH2NH-;
        (4-SMe-phenyl) CH2NH-; (2,3-diCl-phenyl) CH2NH-;
        (2,4-diCl-phenyl)CH2NH-; (2,5-diCl-phenyl)CH2NH-;
        (2,6-diCl-phenyl)CH2NH-; (3,4-diCl-phenyl)CH2NH-;
        (3,5-diCl-phenyl)CH2NH-; (2,3-diF-phenyl)CH2NH-;
10
        (2,4-diF-phenyl) CH2NH-; (2,5-diF-phenyl) CH2NH-;
        (2,6-diF-phenyl)CH2NH-; (3,4-diF-phenyl)CH2NH-;
        (3,5-diF-phenyl)CH2NH-; (2,3-diCH3-phenyl)CH2NH-;
        (2,4-diCH3-phenyl)CH2NH-; (2,5-diCH3-phenyl)CH2NH-;
        (2,6-diCH3-phenyl)CH2NH-; (3,4-diCH3-phenyl)CH2NH-;
15
        (3,5-diCH3-phenyl)CH2NH-; (2,3-diCF3-phenyl)CH2NH-;
        (2,4-diCF3-phenyl)CH2NH-; (2,5-diCF3-phenyl)CH2NH-;
        (2,6-diCF3-phenyl)CH2NH-; (3,4-diCF3-phenyl)CH2NH-;
        (3,5-diCF3-phenyl)CH2NH-; (2,3-diOMe-phenyl)CH2NH-;
        (2,4-diOMe-phenyl)CH2NH-; (2,5-diOMe-phenyl)CH2NH-;
20
       (2,6-diOMe-phenyl)CH2NH-; (3,4-diOMe-phenyl)CH2NH-;
       (3,5-diOMe-phenyl)CH2NH-; (2-F-3-Cl-phenyl)CH2NH-;
       (2-F-4-Cl-phenyl)CH2NH-; (2-F-5-Cl-phenyl)CH2NH-;
       (2-F-6-C1-pheny1)CH2NH-; (2-F-3-CH3-pheny1)CH2NH-;
       (2-F-4-CH3-phenyl)CH2NH-; (2-F-5-CH3-phenyl)CH2NH-;
25
       (2-F-6-CH3-phenyl) CH2NH-; (2-F-3-CF3-phenyl) CH2NH-;
       (2-F-4-CF3-phenyl)CH2NH-; (2-F-5-CF3-phenyl)CH2NH-;
       (2-F-6-CF3-phenyl)CH2NH-; (2-F-3-OMe-phenyl)CH2NH-;
       (2-F-4-OMe-phenyl)CH2NH-; (2-F-5-OMe-phenyl)CH2NH-;
       (2-F-6-OMe-phenyl)CH2NH-; (2-Cl-3-F-phenyl)CH2NH-;
30
       (2-Cl-4-F-phenyl)CH2NH-; (2-Cl-5-F-phenyl)CH2NH-;
       (2-Cl-6-F-phenyl)CH2NH-; (2-Cl-3-CH3-phenyl)CH2NH-;
       (2-C1-4-CH3-phenyl)CH2NH-; (2-C1-5-CH3-phenyl)CH2NH-;
       (2-C1-6-CH3-phenyl)CH2NH-; (2-C1-3-CF3-phenyl)CH2NH-;
       (2-C1-4-CF3-phenyl)CH2NH-; (2-C1-5-CF3-phenyl)CH2NH-;
35
       (2-Cl-6-CF_3-phenyl)CH_2NH-; (2-Cl-3-OMe-phenyl)CH_2NH-;
       (2-Cl-4-OMe-phenyl) CH<sub>2</sub>NH-; (2-Cl-5-OMe-phenyl) CH<sub>2</sub>NH-;
       (2-Cl-6-OMe-phenyl) CH2NH-; (2-CH3-3-F-phenyl) CH2NH-;
```

```
(2-CH<sub>3</sub>-4-F-phenyl)CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-5-F-phenyl)CH<sub>2</sub>NH-;
               (2-CH3-6-F-phenyl) CH2NH-; (2-CH3-3-Cl-phenyl) CH2NH-;
               (2-CH<sub>3</sub>-4-Cl-phenyl)CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-5-Cl-phenyl)CH<sub>2</sub>NH-;
               (2-CH<sub>3</sub>-6-Cl-phenyl) CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-3-CF<sub>3</sub>-phenyl) CH<sub>2</sub>NH-;
      5
               (2-CH<sub>3</sub>-4-CF<sub>3</sub>-phenyl)CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-5-CF<sub>3</sub>-phenyl)CH<sub>2</sub>NH-;
               (2-CH_3-6-CF_3-pheny1) CH_2NH-; (2-CH_3-3-OMe-pheny1) CH_2NH-;
               (2-CH_3-4-OMe-pheny1) CH_2NH-; (2-CH_3-5-OMe-pheny1) CH_2NH-;
               (2-CH<sub>3</sub>-6-OMe-phenyl) CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-3-F-phenyl) CH<sub>2</sub>NH-;
               (2-CF_3-4-F-phenyl) CH_2NH-; (2-CF_3-5-F-phenyl) CH_2NH-;
     10
               (2-CF_3-6-F-phenyl)CH_2NH-; (2-CF_3-3-Cl-phenyl)CH_2NH-;
               (2-CF<sub>3</sub>-4-Cl-phenyl)CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-5-Cl-phenyl)CH<sub>2</sub>NH-;
               (2-CF_3-6-C1-pheny1)CH_2NH-; (2-CF_3-3-CH_3-pheny1)CH_2NH-;
į.
               (2-CF_3-4-CH_3-phenyl)CH_2NH-; (2-CH_3-5-CF_3-phenyl)CH_2NH-;
£ 19
State of the State of
              (2-CF_3-6-CH_3-pheny1)CH_2NH-; (2-CF_3-3-OMe-pheny1)CH_2NH-;
              (2-CF<sub>3</sub>-4-OMe-phenyl) CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-5-OMe-phenyl) CH<sub>2</sub>NH-;
    15
13
              (2-CF<sub>3</sub>-6-OMe-phenyl)CH<sub>2</sub>NH-; (2-OMe-3-F-phenyl)CH<sub>2</sub>NH-;
10
              (2-OMe-4-F-phenyl)CH2NH-; (2-OMe-5-F-phenyl)CH2NH-;
W
10 marie
              (2-OMe-6-F-phenyl)CH2NH-; (2-OMe-3-Cl-phenyl)CH2NH-;
              (2-OMe-4-Cl-phenyl) CH2NH-; (2-OMe-5-Cl-phenyl) CH2NH-;
Part of
              (2-OMe-6-Cl-phenyl) CH2NH-; (2-OMe-4-CN-phenyl) CH2NH-;
    20
43
              (2-OMe-4-CHO-phenyl) CH<sub>2</sub>NH-; (2-OMe-3-CH<sub>3</sub>-phenyl) CH<sub>2</sub>NH-;
0
              (2-OMe-4-CH_3-phenyl) CH_2NH-; (2-OMe-5-CH_3-phenyl) CH_2NH-;
              (2-OMe-6-CH_3-phenyl)CH_2NH-; (2-OMe-3-CF_3-phenyl)CH_2NH-;
              (2-OMe-4-CF3-phenyl) CH2NH-; (2-OMe-5-CF3-phenyl) CH2NH-;
    25
              (2-OMe-6-CF3-phenyl) CH2NH-; (2-acetyl-4-Cl-phenyl) CH2NH-;
              (2-acetyl-4-Me-phenyl)CH2NH-;
              (2-acetyl-4-MeO-phenyl) CH2NH-;
              (2-CH3CH(OH)-4-Cl-phenyl)CH2NH-;
              (2-CH3CH(OH)-4-Me-phenyl)CH2NH-;
    30
              (2-CH3CH(OH)-4-MeO-phenyl)CH2NH-;
              (3-CF<sub>3</sub>-4-Cl-phenyl)CH<sub>2</sub>NH-; (3-F-4-CHO-phenyl)CH<sub>2</sub>NH-;
              (3-CH_3-4-CN-phenyl)CH_2NH-; (3-CH_3-4-MeO-phenyl)CH_2NH-;
              (3-CH<sub>3</sub>-4-Cl-phenyl)CH<sub>2</sub>NH-; (3-CH<sub>3</sub>-4-F-phenyl)CH<sub>2</sub>NH-;
    35
              (4-F-3-CF<sub>3</sub>-phenyl)CH<sub>2</sub>NH-; (3-CH<sub>3</sub>-4-CO<sub>2</sub>Me-phenyl)CH<sub>2</sub>NH-;
              (3-CF3-4-C(0)CH3-phenyl)CH2NH-;
              (3-CHO-4-OMe-phenyl) CH2NH-;
```

5

```
(2,3,5-triCl-phenyl)CH<sub>2</sub>NH-;
(2,4,5-triF-phenyl)CH<sub>2</sub>NH-;
(2,6-diCl-3-Me-phenyl)CH<sub>2</sub>NH-;
(3,5-diMe-4-MeO-phenyl)CH<sub>2</sub>NH-; and
(2-F-3-Cl-6-CF<sub>3</sub>-phenyl)CH<sub>2</sub>NH-;
```

provided that two of R⁷, R⁸, and R⁹, are independently selected from hydrogen, fluoro, chloro, bromo, cyano, 10 methyl, ethyl, propyl, isopropyl, butyl, t-butyl, nitro, trifluoromethyl, methoxy, ethoxy, isopropoxy, and trifluoromethoxy.

8. A compound of Claim 7 of Formula (II)

wherein:

20 b is a single bond, wherein the bridge hydrogens are in a cis or trans position;

R1 is selected from

hydrogen, methyl, ethyl, n-propyl, n-butyl, s-butyl,
t-butyl, n-pentyl, n-hexyl, 2-propyl, 2-butyl, 2-pentyl,
2-hexyl, 2-methylpropyl, 2-methylbutyl, 2-methylpentyl,
2-ethylbutyl, 3-methylpentyl, 3-methylbutyl,
4-methylpentyl, 2-fluoroethyl, 2,2-difluoroethyl,
2,2-trifluoroethyl, 2-propenyl, 2-methyl-2-propenyl,
trans-2-butenyl, 3-methyl-2-butenyl, 3-butenyl,
trans-2-pentenyl, cis-2-pentenyl, 4-pentenyl,
4-methyl-3-pentenyl, 3,3-dichloro-2-propenyl,

```
trans-3-phenyl-2-propenyl, cyclopropyl, cyclobutyl,
        cyclopentyl, cyclohexyl, cyclopropylmethyl,
        cyclobutylmethyl, cyclopentylmethyl, cyclohexylmethyl,
        -CH=CH_2, -CH_2-CH=CH_2, -CH=CH-CH_3, -C\equiv CH, -C\equiv C-CH_3,
 5
        and -CH2-C=CH;
     R6a is H:
     R6b is H;
10
     alternatively, R^{6a} and R^{6b} are taken together to form =0;
     {\ensuremath{\mathsf{R}}}^7 and {\ensuremath{\mathsf{R}}}^9, at each occurrence, are independently selected
           from hydrogen, fluoro, methyl, trifluoromethyl, and
           methoxy;
     R8 is selected from
        hydrogen, fluoro, chloro, bromo, cyano, methyl, ethyl,
        propyl, isopropyl, butyl, t-butyl, nitro,
        trifluoromethyl, methoxy, ethoxy, isopropoxy,
        trifluoromethoxy, phenyl;
        2-Cl-phenyl; 2-F-phenyl; 2-Br-phenyl; 2-CN-phenyl;
        2-Me-phenyl; 2-CF3-phenyl; 2-MeO-phenyl; 2-CF3O-phenyl;
25
       2-NO<sub>2</sub>-phenyl; 2-MeS-phenyl; 2-CHO-phenyl; 2-HOCH<sub>2</sub>-
       phenyl;
       3-Cl-phenyl; 3-F-phenyl; 3-Br-phenyl; 3-CN-phenyl;
       3-Me-phenyl; 3-Et-phenyl; 3-n-Pr-phenyl; 3-isoPr-phenyl;
30
       3-n-Bu-phenyl; 3-CF3-phenyl; 3-MeO-phenyl; 3-MeS-phenyl;
       3-isopropoxyphenyl; 3-CF3O-phenyl; 3-NO2-phenyl;
       3-CHO-phenyl; 3-HOCH2-phenyl; 3-MeOCH2-phenyl;
       3-MeaNCHa-phenvl:
       4-Cl-phenyl; 4-F-phenyl; 4-Br-phenyl; 4-CN-phenyl;
35
       4-Me-phenyl; 4-Et-phenyl; 4-n-Pr-phenyl; 4-iso-Pr-
       phenyl;
```

```
4-n-Bu-phenyl; 4-CF3-phenyl; 4-MeO-phenyl;
        4-isopropoxyphenyl; 4-CF3O-phenyl; 4-MeS-phenyl;
        4-acetylphenyl; 3-acetamidophenyl; 4-pyridyl;
 5
        2-furanyl; 2-thiophenyl; 2-naphthyl; 1-pyrrolidinyl,
        2,3-diCl-phenyl; 2,3-diF-phenyl; 2,3-diMe-phenyl;
       2,3-diCF3-phenyl; 2,3-diMeO-phenyl; 2,3-diCF3O-phenyl;
10
       2,4-diCl-phenyl; 2,4-diF-phenyl; 2,4-diMe-phenyl;
       2,4-diCF3-phenyl; 2,4-diMeO-phenyl; 2,4-diCF3O-phenyl;
       2,5-diCl-phenyl; 2,5-diF-phenyl; 2,5-diMe-phenyl;
       2,5-diCF3-phenyl; 2,5-diMeO-phenyl; 2,5-diCF3O-phenyl;
15
       2,6-diCl-phenyl; 2,6-diF-phenyl; 2,6-diMe-phenyl;
       2,6-diCF3-phenyl; 2,6-diMeO-phenyl; 2,6-diCF3O-phenyl;
       3,4-diCl-phenyl; 3,4-diF-phenyl; 3,4-diMe-phenyl;
20
       3,4-diCF3-phenyl; 3,4-diMeO-phenyl; 3,4-diCF3O-phenyl;
       2,4,6-triCl-phenyl; 2,4,6-triF-phenyl;
       2,4,6-triMe-phenyl; 2,4,6-triCF3-phenyl;
       2,4,6-triMeO-phenyl; 2,4,6-triCF3O-phenyl;
25
       2,4,5-triMe-phenyl; 2,3,4-triF-phenyl;
       2-Me-4-MeO-5-F-phenyl; 2,6-diCl-4-MeO-phenyl;
       2,4-diMeO-6-F-phenyl; 2,6-diF-4-Cl-phenyl;
       2,3,4,6-tetraF-phenyl; 2,3,4,5,6-pentaF-phenyl;
30
       2-Cl-4-F-phenyl; 2-Cl-6-F-phenyl; 2-Cl-3-Me-phenyl;
       2-Cl-4-MeO-phenyl; 2-Cl-4-EtO-phenyl;
       2-Cl-4-iPrO-phenyl; 2-Cl-4-CF3-phenyl;
       2-Cl-4-CF3O-phenyl; 2-Cl-4-(CHF2)O-phenyl;
       2-F-3-Cl-phenyl; 2-F-4-MeO-phenyl; 2-F-5-Me-phenyl;
35
       2-Me-3-Cl-phenyl; 2-Me-3-CN-phenyl; 2-Me-4-Cl-phenyl;
       2-Me-4-F-phenyl; 2-Me-4-CN-phenyl; 2-Me-4-MeO-phenyl;
```

```
2-Me-4-EtO-phenyl; 2-Me-4-MeS-phenyl;
        2-Me-4-H2NCO-phenyl; 2-Me-4-MeOC(=O)-phenyl;
        2-Me-4-CH<sub>3</sub>C(=0)-phenyl; 2-Me-5-F-phenyl;
        2-Et-4-MeO-phenvl: 2-MeO-5-F-phenvl:
 5
        2-MeO-4-isopropyl-phenyl; 2-CF3-4-Cl-phenyl;
        2-CF3-4-F-phenyl; 2-CF3-4-MeO-phenyl;
        2-CF3-4-EtO-phenyl; 2-CF3-4-iPrO-phenyl;
        2-CF3-4-CN-phenyl; 2-CF3-6-F-phenyl;
        2-CHO-4-MeO-phenyl; 2-MeOC(=0)-3-MeO-phenyl;
10
        2-CH<sub>3</sub>CH(OH)-4-MeO-phenyl; 2-CH<sub>3</sub>CH(OH)-4-F-phenyl;
        2-CH3CH(OH)-4-Cl-phenyl; 2-CH3CH(OH)-4-Me-phenyl;
        2-CH<sub>3</sub>CH(OMe)-4-MeO-phenyl; 2-CH<sub>3</sub>C(=O)-4-MeO-phenyl;
        2-CH3C(=0)-4-F-phenyl; 2-CH3C(=0)-4-Cl-phenyl;
        2-CH<sub>3</sub>C(=0)-4-Me-phenyl; 2-H<sub>2</sub>C(OH)-4-MeO-phenyl;
        2-H<sub>2</sub>C(OMe)-4-MeO-phenyl; 2-H<sub>3</sub>CCH<sub>2</sub>CH(OH)-4-MeO-phenyl;
15
        2-H3CCH2C(=0)-4-MeO-phenyl; 2-CH3CO2CH2CH2-4-MeO-phenyl;
        (Z) -2-HOCH2CH=CH-4-MeO-phenyl;
        (E) -2-HOCH2CH=CH-4-MeO-phenvl;
        (Z) -2-CH3CO2CH=CH-4-MeO-phenyl;
20
        (E) -2-CH3CO2CH=CH-4-MeO-phenyl;
        2-CH3OCH2CH2-4-MeO-phenyl;
        3-CN-4-F-phenyl; 3-H2NCO-4-F-phenyl;
        (2-Cl-phenyl)-CH=CH-; (3-Cl-phenyl)-CH=CH-;
25
        (2,6-diF-phenyl)-CH=CH-; phenyl-CH=CH-;
        (2-Me-4-MeO-phenvl)-CH=CH-:
       cyclohexyl; cyclopentyl; cyclohexylmethyl; benzyl;
       2-F-benzyl; 3-F-benzyl; 4-F-benzyl; 3-MeO-benzyl;
30
       3-OH-benzyl; 2-MeO-benzyl; 2-OH-benzyl;
       tetrahydroguinolin-1-vl:
       tetrahydroindolin-1-vl;
       tetrahydroisoindolin-1-yl;
35
       phenyl-S-; phenyl-NH-; pyrid-3-yl-NH-;
       (4-Me-pyrid-3-yl)-NH-; (4-Cl-pyrid-3-yl)-NH-;
       (1-naphthyl)-NH-; (2-naphthyl)-NH-;
```

```
(2-Me-naphth-1-yl)-NH-; (4-Me-naphth-1-yl)-NH-;
        (3-quinolinvl)-NH-:
        (2-[1,1'-biphenyl])-NH-; (3-[1,1'-biphenyl])-NH-;
 5
       (4-[1,1'-biphenyl])-NH-; (2-F-phenyl)-NH-;
       (2-Cl-phenyl)-NH-; (2-CF3-phenyl)-NH-;
       (2-CH3-phenyl)-NH-; (2-OMe-phenyl)-NH-;
       (2-CN-phenyl)-NH-; (2-OCF3-phenyl)-NH-;
       (2-SMe-phenyl)-NH-; (3-F-phenyl)-NH-;
       (3-Cl-phenyl)-NH-; (3-CF3-phenyl)-NH-;
10
       (3-CH3-phenyl)-NH-; (3-OMe-phenyl)-NH-;
       (3-CN-phenyl)-NH-; (3-OCF3-phenyl)-NH-;
       (3-SMe-phenyl)-NH-; (4-F-phenyl)-NH-;
       (4-Cl-phenyl)-NH-; (4-CF3-phenyl)-NH-;
15
       (4-CH3-phenyl)-NH-; (4-OMe-phenyl)-NH-;
       (4-CN-phenyl)-NH-; (4-OCF3-phenyl)-NH-;
       (4-SMe-phenyl)-NH-; (2,3-diCl-phenyl)-NH-;
       (2,4-diCl-phenyl)-NH-; (2,5-diCl-phenyl)-NH-;
       (2,6-diCl-phenyl)-NH-; (3,4-diCl-phenyl)-NH-;
       (3,5-diCl-phenyl)-NH-; (2,3-diF-phenyl)-NH-;
20
       (2,4-diF-phenyl)-NH-; (2,5-diF-phenyl)-NH-;
       (2,6-diF-phenyl)-NH-; (3,4-diF-phenyl)-NH-;
       (3,5-diF-phenyl)-NH-; (2,3-diCH3-phenyl)-NH-;
       (2,4-diCH3-phenyl)-NH-; (2,5-diCH3-phenyl)-NH-;
25
       (2,6-diCH3-phenyl)-NH-; (3,4-diCH3-phenyl)-NH-;
       (3,5-diCH3-phenyl)-NH-; (2,3-diCF3-phenyl)-NH-;
       (2,4-diCF3-phenyl)-NH-; (2,5-diCF3-phenyl)-NH-;
       (2,6-diCF3-phenyl)-NH-; (3,4-diCF3-phenyl)-NH-;
       (3,5-diCF3-phenyl)-NH-; (2,3-diOMe-phenyl)-NH-;
30
       (2,4-diOMe-phenyl)-NH-; (2,5-diOMe-phenyl)-NH-;
       (2,6-diOMe-phenyl)-NH-; (3,4-diOMe-phenyl)-NH-;
       (3,5-diOMe-phenyl)-NH-; (2-F-3-Cl-phenyl)-NH-;
       (2-F-4-Cl-phenyl)-NH-; (2-F-5-Cl-phenyl)-NH-;
       (2-F-6-Cl-phenyl)-NH-; (2-F-3-CH3-phenyl)-NH-;
35
       (2-F-4-CH3-phenyl)-NH-; (2-F-5-CH3-phenyl)-NH-;
       (2-F-6-CH3-phenyl)-NH-; (2-F-3-CF3-phenyl)-NH-;
       (2-F-4-CF3-phenyl)-NH-; (2-F-5-CF3-phenyl)-NH-;
```

£ 44

```
(2-F-6-CF3-phenyl)-NH-; (2-F-3-OMe-phenyl)-NH-;
         (2-F-4-OMe-phenyl) -NH-; (2-F-5-OMe-phenyl) -NH-;
         (2-F-6-OMe-phenyl)-NH-; (2-Cl-3-F-phenyl)-NH-;
         (2-Cl-4-F-phenyl)-NH-; (2-Cl-5-F-phenyl)-NH-;
 5
         (2-Cl-6-F-phenyl)-NH-; (2-Cl-3-CH3-phenyl)-NH-;
         (2-Cl-4-CH3-phenyl)-NH-; (2-Cl-5-CH3-phenyl)-NH-;
         (2-Cl-6-CH3-phenyl)-NH-; (2-Cl-3-CF3-phenyl)-NH-;
         (2-Cl-4-CF3-phenyl)-NH-; (2-Cl-5-CF3-phenyl)-NH-;
         (2-Cl-6-CF3-phenyl)-NH-; (2-Cl-3-OMe-phenyl)-NH-;
         (2-Cl-4-OMe-phenyl)-NH-; (2-Cl-5-OMe-phenyl)-NH-;
10
         (2-Cl-6-OMe-phenyl)-NH-; (2-CH3-3-F-phenyl)-NH-;
         (2-CH<sub>3</sub>-4-F-phenyl)-NH-; (2-CH<sub>3</sub>-5-F-phenyl)-NH-;
         (2-CH3-6-F-phenyl)-NH-; (2-CH3-3-Cl-phenyl)-NH-;
         (2-CH<sub>3</sub>-4-Cl-phenyl)-NH-; (2-CH<sub>3</sub>-5-Cl-phenyl)-NH-;
15
         (2-CH<sub>3</sub>-6-Cl-phenyl)-NH-; (2-CH<sub>3</sub>-3-CF<sub>3</sub>-phenyl)-NH-;
         (2-CH3-4-CF3-phenyl)-NH-; (2-CH3-5-CF3-phenyl)-NH-;
        (2-CH3-6-CF3-phenyl)-NH-; (2-CH3-3-OMe-phenyl)-NH-;
        (2-CH<sub>3</sub>-4-OMe-phenyl)-NH-; (2-CH<sub>3</sub>-5-OMe-phenyl)-NH-;
        (2-CH<sub>3</sub>-6-OMe-phenyl)-NH-; (2-CF<sub>3</sub>-3-F-phenyl)-NH-;
20
        (2-CF3-4-F-phenyl)-NH-; (2-CF3-5-F-phenyl)-NH-;
        (2-CF<sub>3</sub>-6-F-phenyl)-NH-; (2-CF<sub>3</sub>-3-Cl-phenyl)-NH-;
        (2-CF<sub>3</sub>-4-Cl-phenyl)-NH-; (2-CF<sub>3</sub>-5-Cl-phenyl)-NH-;
        (2-CF3-6-Cl-phenyl)-NH-; (2-CF3-3-CH3-phenyl)-NH-;
        (2-CF<sub>3</sub>-4-CH<sub>3</sub>-phenyl)-NH-; (2-CH<sub>3</sub>-5-CF<sub>3</sub>-phenyl)-NH-;
25
        (2-CF<sub>3</sub>-6-CH<sub>3</sub>-phenyl)-NH-; (2-CF<sub>3</sub>-3-OMe-phenyl)-NH-;
        (2-CF<sub>3</sub>-4-OMe-phenyl)-NH-; (2-CF<sub>3</sub>-5-OMe-phenyl)-NH-;
        (2-CF<sub>3</sub>-6-OMe-phenyl)-NH-; (2-OMe-3-F-phenyl)-NH-;
        (2-OMe-4-F-phenyl)-NH-; (2-OMe-5-F-phenyl)-NH-;
        (2-OMe-6-F-phenyl)-NH-; (2-OMe-3-Cl-phenyl)-NH-;
30
        (2-OMe-4-Cl-phenyl)-NH-; (2-OMe-5-Cl-phenyl)-NH-;
        (2-OMe-6-Cl-phenyl)-NH-; (2-OMe-4-CN-phenyl)-NH-;
        (2-OMe-4-CHO-phenyl)-NH-; (2-OMe-3-CH3-phenyl)-NH-;
        (2-OMe-4-CH3-phenyl)-NH-; (2-OMe-5-CH3-phenyl)-NH-;
        (2-OMe-6-CH3-phenyl)-NH-; (2-OMe-3-CF3-phenyl)-NH-;
35
        (2-OMe-4-CF3-phenyl)-NH-; (2-OMe-5-CF3-phenyl)-NH-;
        (2-OMe-6-CF3-phenyl)-NH-; (2-acetyl-4-Cl-phenyl)-NH-;
        (2-acetyl-4-Me-phenyl)-NH-; (2-acetyl-4-MeO-phenyl)-NH-;
```

```
(2-CH3CH(OH)-4-Cl-phenvl)-NH-:
        (2-CH3CH(OH)-4-Me-phenvl)-NH-;
        (2-CH3CH(OH)-4-MeO-phenyl)-NH-;
 5
        (3-CF<sub>3</sub>-4-Cl-phenyl)-NH-; (3-F-4-CHO-phenyl)-NH-;
        (3-CH<sub>3</sub>-4-CN-phenyl)-NH-; (3-CH<sub>3</sub>-4-MeO-phenyl)-NH-;
        (3-CH3-4-Cl-phenyl)-NH-; (3-CH3-4-F-phenyl)-NH-;
        (3-F-5-CF3-phenyl)-NH-;
10
        (3-CH<sub>3</sub>-4-CO<sub>2</sub>Me-phenyl)NH-; (3-CF<sub>3</sub>-4-C(O)CH<sub>3</sub>-phenyl)NH-;
        (3-CHO-4-OMe-phenyl)-NH-; (4-F-3-CF3-phenyl)-NH-;
        (2,3,5-triCl-phenyl)-NH-; (2,4,5-triF-phenyl)-NH-;
        (2,6-diCl-3-Me-phenyl)-NH-; (3,5-diMe-4-MeO-phenyl)-NH-;
15
        (2-F-3-C1-6-CF3-phenyl)-NH-;
        benzyl-NH-; (3-quinolinyl)CH2NH-; (2-F-phenyl)CH2NH-;
        (2-Cl-phenyl) CH2NH-; (2-CF3-phenyl) CH2NH-;
        (2-CH3-phenyl) CH2NH-; (2-OMe-phenyl) CH2NH-;
20
        (2-CN-phenyl) CH2NH-; (2-OCF3-phenyl) CH2NH-;
        (2-SMe-phenyl) CH2NH-; (3-F-phenyl) CH2NH-;
        (3-Cl-phenyl) CH2NH-; (3-CF3-phenyl) CH2NH-;
        (3-CH3-phenyl) CH2NH-; (3-OMe-phenyl) CH2NH-;
        (3-CN-phenyl) CH2NH-; (3-OCF3-phenyl) CH2NH-;
        (3-SMe-phenyl)CH2NH-; (4-F-phenyl)CH2NH-;
25
        (4-Cl-phenyl) CH2NH-; (4-CF3-phenyl) CH2NH-;
        (4-CH3-phenyl) CH2NH-; (4-OMe-phenyl) CH2NH-;
        (4-CN-phenyl) CH2NH-; (4-OCF3-phenyl) CH2NH-;
        (4-SMe-phenyl) CH2NH-; (2,3-diCl-phenyl) CH2NH-;
30
        (2,4-diCl-phenyl) CH2NH-; (2,5-diCl-phenyl) CH2NH-;
        (2,6-diCl-phenyl)CH2NH-; (3,4-diCl-phenyl)CH2NH-;
        (3,5-diCl-phenyl)CH2NH-; (2,3-diF-phenyl)CH2NH-;
        (2,4-diF-phenyl)CH2NH-; (2,5-diF-phenyl)CH2NH-;
        (2,6-diF-phenyl)CH2NH-; (3,4-diF-phenyl)CH2NH-;
35
        (3,5-diF-phenyl)CH2NH-; (2,3-diCH3-phenyl)CH2NH-;
        (2,4-diCH3-phenyl)CH2NH-; (2,5-diCH3-phenyl)CH2NH-;
        (2,6-diCH3-phenyl)CH2NH-; (3,4-diCH3-phenyl)CH2NH-;
```

```
(3,5-diCH3-phenyl)CH2NH-; (2,3-diCF3-phenyl)CH2NH-;
         (2,4-diCF3-phenyl)CH2NH-; (2,5-diCF3-phenyl)CH2NH-;
         (2,6-diCF3-phenyl)CH2NH-; (3,4-diCF3-phenyl)CH2NH-;
         (3,5-diCF3-phenyl)CH2NH-; (2,3-diOMe-phenyl)CH2NH-;
 5
         (2,4-diOMe-phenyl)CH2NH-; (2,5-diOMe-phenyl)CH2NH-;
         (2,6-diOMe-phenyl)CH2NH-; (3,4-diOMe-phenyl)CH2NH-;
         (3,5-diOMe-phenyl)CH2NH-; (2-F-3-Cl-phenyl)CH2NH-;
         (2-F-4-Cl-phenyl)CH2NH-; (2-F-5-Cl-phenyl)CH2NH-;
         (2-F-6-Cl-phenyl)CH2NH-; (2-F-3-CH3-phenyl)CH2NH-;
10
         (2-F-4-CH3-phenyl) CH2NH-; (2-F-5-CH3-phenyl) CH2NH-;
         (2-F-6-CH3-phenyl) CH2NH-; (2-F-3-CF3-phenyl) CH2NH-;
         (2-F-4-CF3-phenyl)CH2NH-; (2-F-5-CF3-phenyl)CH2NH-;
         (2-F-6-CF3-phenyl)CH2NH-; (2-F-3-OMe-phenyl)CH2NH-;
         (2-F-4-OMe-phenyl)CH2NH-; (2-F-5-OMe-phenyl)CH2NH-;
15
         (2-F-6-OMe-phenyl) CH2NH-; (2-Cl-3-F-phenyl) CH2NH-;
         (2-Cl-4-F-phenyl) CH2NH-; (2-Cl-5-F-phenyl) CH2NH-;
         (2-Cl-6-F-phenyl) CH2NH-; (2-Cl-3-CH3-phenyl) CH2NH-;
         (2-Cl-4-CH<sub>3</sub>-phenyl)CH<sub>2</sub>NH-; (2-Cl-5-CH<sub>3</sub>-phenyl)CH<sub>2</sub>NH-;
         (2-Cl-6-CH3-phenyl) CH2NH-; (2-Cl-3-CF3-phenyl) CH2NH-;
20
         (2-Cl-4-CF3-phenyl) CH2NH-; (2-Cl-5-CF3-phenyl) CH2NH-;
         (2-Cl-6-CF3-phenyl) CH2NH-; (2-Cl-3-OMe-phenyl) CH2NH-;
         (2-Cl-4-OMe-phenyl) CH2NH-; (2-Cl-5-OMe-phenyl) CH2NH-;
         (2-Cl-6-OMe-phenyl) CH2NH-; (2-CH3-3-F-phenyl) CH2NH-;
         (2-CH<sub>3</sub>-4-F-phenyl) CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-5-F-phenyl) CH<sub>2</sub>NH-;
25
         (2-CH3-6-F-phenyl)CH2NH-; (2-CH3-3-Cl-phenyl)CH2NH-;
         (2-CH3-4-Cl-phenyl) CH2NH-; (2-CH3-5-Cl-phenyl) CH2NH-;
         (2-CH3-6-Cl-phenyl) CH2NH-; (2-CH3-3-CF3-phenyl) CH2NH-;
         (2-CH2-4-CF2-phenvl) CH2NH-; (2-CH2-5-CF3-phenyl) CH2NH-;
         (2-CH<sub>3</sub>-6-CF<sub>3</sub>-phenyl)CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-3-OMe-phenyl)CH<sub>2</sub>NH-;
30
        (2-CH<sub>3</sub>-4-OMe-phenyl)CH<sub>2</sub>NH-; (2-CH<sub>3</sub>-5-OMe-phenyl)CH<sub>2</sub>NH-;
        (2-CH_3-6-OMe-phenyl)CH_2NH-; (2-CF_3-3-F-phenyl)CH_2NH-;
        (2-CF<sub>3</sub>-4-F-phenyl) CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-5-F-phenyl) CH<sub>2</sub>NH-;
        (2-CF_3-6-F-pheny1)CH_2NH-; (2-CF_3-3-C1-pheny1)CH_2NH-;
        (2-CF<sub>3</sub>-4-Cl-phenyl)CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-5-Cl-phenyl)CH<sub>2</sub>NH-;
35
        (2-CF<sub>3</sub>-6-Cl-phenyl)CH<sub>2</sub>NH-; (2-CF<sub>3</sub>-3-CH<sub>3</sub>-phenyl)CH<sub>2</sub>NH-;
        (2-CF3-4-CH3-phenyl)CH2NH-; (2-CH3-5-CF3-phenyl)CH2NH-;
        (2-CF3-6-CH3-phenyl) CH2NH-; (2-CF3-3-OMe-phenyl) CH2NH-;
```

```
(2-CF3-4-OMe-phenyl) CH2NH-; (2-CF3-5-OMe-phenyl) CH2NH-;
        (2-CF3-6-OMe-phenyl) CH2NH-; (2-OMe-3-F-phenyl) CH2NH-;
        (2-OMe-4-F-phenyl)CH2NH-; (2-OMe-5-F-phenyl)CH2NH-;
        (2-OMe-6-F-phenyl) CH2NH-; (2-OMe-3-Cl-phenyl) CH2NH-;
        (2-OMe-4-Cl-phenyl) CH2NH-; (2-OMe-5-Cl-phenyl) CH2NH-;
 5
        (2-OMe-6-Cl-phenyl) CH2NH-; (2-OMe-4-CN-phenyl) CH2NH-;
        (2-OMe-4-CHO-phenyl) CH2NH-; (2-OMe-3-CH3-phenyl) CH2NH-;
        (2-OMe-4-CH3-phenyl) CH2NH-; (2-OMe-5-CH3-phenyl) CH2NH-;
        (2-OMe-6-CH3-phenyl) CH2NH-; (2-OMe-3-CF3-phenyl) CH2NH-;
10
        (2-OMe-4-CF3-phenyl) CH2NH-; (2-OMe-5-CF3-phenyl) CH2NH-;
        (2-OMe-6-CF3-phenyl) CH2NH-; (2-acetyl-4-Cl-phenyl) CH2NH-;
        (2-acetyl-4-Me-phenyl)CHoNH-;
        (2-acetyl-4-MeO-phenyl)CH2NH-;
        (2-CH3CH(OH)-4-Cl-phenyl)CH2NH-;
15
        (2-CH3CH(OH)-4-Me-phenyl)CH2NH-;
        (2-CH3CH(OH)-4-MeO-phenvl)CH2NH-:
        (3-CF3-4-Cl-phenyl) CH2NH-; (3-F-4-CHO-phenyl) CH2NH-;
        (3-CH<sub>3</sub>-4-CN-phenyl) CH<sub>2</sub>NH-; (3-CH<sub>3</sub>-4-MeO-phenyl) CH<sub>2</sub>NH-;
20
        (3-CH<sub>3</sub>-4-Cl-phenyl) CH<sub>2</sub>NH-; (3-CH<sub>3</sub>-4-F-phenyl) CH<sub>2</sub>NH-;
        (4-F-3-CF_3-phenyl) CH<sub>2</sub>NH-; (3-CH_3-4-CO_2Me-phenyl) CH<sub>2</sub>NH-;
        (3-CF3-4-C(0)CH3-phenyl)CH2NH-;
        (3-CHO-4-OMe-phenyl) CH2NH-;
25
        (2,3,5-triCl-phenyl)CH2NH-;
        (2,4,5-triF-phenyl)CH2NH-;
        (2,6-diC1-3-Me-phenv1)CHoNH-;
        (3,5-diMe-4-MeO-phenyl) CH2NH-; and
        (2-F-3-C1-6-CF3-phenyl)CH2NH-.
30
          A compound of Claim 1, 2, 3, 4, 5, 6, or 7, wherein X
     9.
           is a bond.
```

10. A compound of Claim 1, 2, 3, 4, 5, 6, or 7, wherein X is -O- or -S-.

- 11. A compound of Claim 1, 2, 3, 4, 5, 6, or 7, wherein X is -OCH₂- or -SCH₂-.
- 12. A compound of Claim 1, 2, 3, 4, 5, 6, or 7, wherein X is $-CH_2-$.
 - 13. A compound of Claim 1 wherein:
- 10 X is a bond, $-CH_2-$, -O-, -S-, -S(=O)-, $-S(=O)_2-$, $-NR^{10}-$, $-CH_2CH_2-$, $-CCH_2-$, $-SCH_2-$, $-CH_2O-$, $-CH_2S-$, or $-CH_2NR^{10}-$;
 - R1 is selected from
 - C1-6 alkyl substituted with Z,
 - C_{2-6} alkenyl substituted with Z,
 - C2-6 alkynyl substituted with Z,
 - C3-6 cycloalkyl substituted with Z,
 - aryl substituted with Z,
 - 5-6 membered heterocyclic ring system containing at least one heteroatom selected from the group consisting of N, O, and S, said heterocyclic ring system substituted with Z;
 - C1-6 alkyl substituted with 0-2 R2,
 - C_{2-6} alkenyl substituted with 0-2 \mathbb{R}^2 ,
- 25 C₂₋₆ alkynyl substituted with 0-2 R²,
 - aryl substituted with 0-2 R2, and
 - 5-6 membered heterocyclic ring system containing at least one heteroatom selected from the group consisting of N, O, and S, said heterocyclic ring system substituted with 0-2 R²;
- Z is selected from H,
 - -CH(OH)R2,
 - -C(ethylenedioxy)R2,
- 35 -OR²,

- -SR2,
- $-NR^2R^3$.

```
-C(0)R2,
               -C(0)NR2R3,
               -NR3C(0)R2.
               -C(0)OR2,
     5
               -OC(O)R2,
               -CH (=NR^4) NR^2R^3.
               -NHC (=NR^4) NR^2R^3
               -S(0)R2,
               -S(0)2R2,
    10
               -S(0)_2NR^2R^3, and -NR^3S(0)_2R^2:
         {\ensuremath{R}}^2, at each occurrence, is independently selected from
Li
0
               C_{1-4} alkyl,
13
C2-4 alkenyl,
               C_{2-4} alkynyl,
               C3-6 cycloalkyl,
               aryl substituted with 0-5 R42:
ill sales
               C_{3-10} carbocyclic residue substituted with 0-3 R^{41}, and
               5-10 membered heterocyclic ring system containing from
them today and Book with
   20
                     1-4 heteroatoms selected from the group
                     consisting of N, O, and S substituted with 0-3
                     R41;
         \mathbb{R}^3, at each occurrence, is independently selected from
                H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, and
   25
               C_{1-4} alkoxy;
         alternatively, \mathbb{R}^2 and \mathbb{R}^3 join to form a 5- or 6-membered
               ring optionally substituted with -O- or -N(R^4)-;
   30
        \mathbb{R}^4, at each occurrence, is independently selected from H,
               methyl, ethyl, propyl, and butyl;
        R6a is H or C1-4 alkyl;
   35
        R6b is H;
```

25

30

35

```
alternatively, R6a and R6b are taken together to form =0 or
     =S;
```

R7, R8, and R9, at each occurrence, are independently selected from H, halo, -CF₃, -OCF₃, -OH, -CN, -NO₂, -NR⁴⁶R⁴⁷, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl, C_{1-8} alkoxy, $(C_{1-4}$ haloalkyl)oxy, C_{1-4} alkyl substituted with 0-2 R^{11} . 10 C₃₋₁₀ carbocyclic residue substituted with 0-3 R³³,

arvl substituted with 0-5 R33, 5-10 membered heterocyclic ring system containing from

1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R31:

 OR^{12} , SR^{12} , $NR^{12}R^{13}$, C(O)H, $C(O)R^{12}$, $C(O)NR^{12}R^{13}$, $NR^{14}C(0)R^{12}$, $C(0)OR^{12}$, $OC(0)R^{12}$, $OC(0)OR^{12}$, $CH(=NR^{14})NR^{12}R^{13}$, $NHC(=NR^{14})NR^{12}R^{13}$, $S(O)R^{12}$, $S(O)_2R^{12}$, $S(O)NR^{12}R^{13}$, $S(O)_2NR^{12}R^{13}$, $NR^{14}S(O)R^{12}$, $NR^{14}S(O)_2R^{12}$, $NR^{12}C(0)R^{15}$, $NR^{12}C(0)OR^{15}$, $NR^{12}S(0)_2R^{15}$, and NR12C(O)NHR15:

 R^{10} is selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, and C1-4 alkoxy;

R11 is selected from

H, halo, -CF3, -CN, -NO2,

 C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} haloalkyl, C1-8 alkoxy, C3-10 cycloalkyl,

C₃₋₁₀ carbocyclic residue substituted with 0-3 R³³, aryl substituted with 0-5 R33,

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R31:

35

5

10

```
OR<sup>12</sup>, SR<sup>12</sup>, NR<sup>12</sup>R<sup>13</sup>, C(O)H, C(O)R<sup>12</sup>, C(O)NR<sup>12</sup>R<sup>13</sup>, NR^{14}C(O)R^{12}, C(O)OR^{12}, OC(O)R^{12}, OC(O)R^{12}, OC(O)R^{12}, \\ CH(=NR^{14})NR^{12}R^{13}, NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, \\ S(O)_2R^{12}, S(O)NR^{12}R^{13}, S(O)_2NR^{12}R^{13}, NR^{14}S(O)R^{12}, \\ and NR^{14}S(O)_2R^{12};
```

 $R^{12},$ at each occurrence, is independently selected from C_{1-4} alkyl, $C_{2-4} \text{ alkenyl,} \\ C_{2-4} \text{ alkynyl,}$

C₃₋₆ cycloalkyl,

phenyl substituted with 0-5 R^{33} ;

 C_{3-10} carbocyclic residue substituted with 0-3 R^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 R^{31} ;

- R^{13} , at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
- alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or -N(R^{14})-;
- 25 R^{14} , at each occurrence, is independently selected from H and C_{1-4} alkyl;
 - R^{31} , at each occurrence, is independently selected from H, OH, halo, CF₃, SO_2R^{45} , $NR^{46}R^{47}$, methyl, ethyl, and propyl;
 - $R^{33},$ at each occurrence, is independently selected from H, OH, halo, CN, NO $_2$, CF $_3$, SO $_2R^{45}$, NR 4 GR 4 7, C1-3 alkyl, C2-3 alkenyl, C2-3 alkynyl, C3-5 cycloalkyl, C1-3 haloalkyl, C1-3 haloalkyl-oxy-, C1-3 alkyloxy-, C1-3 alkylthio-, C1-3 alkyl-C(=0)-, and C1-3 alkyl-C(=0)NH-:

5

 R^{41} , at each occurrence, is independently selected from H, CF3, halo, OH, CO_2H , SO_2R^{45} , $NR^{46}R^{47}$, NO_2 , CN, =0, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl C_{1-4} alkyl substituted with 0-1 R^{43} , aryl substituted with 0-3 R^{42} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group

1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3

R⁴⁴;

 $R^{42},$ at each occurrence, is independently selected from H, CF3, halo, OH, CO2H, SO2R^45, SR^45, NR^46R^47, OR^48, NO2, CN, CH(=NH)NH2, NHC(=NH)NH2,

 $\text{C}_{2\text{-}6}$ alkenyl, $\text{C}_{2\text{-}6}$ alkynyl, $\text{C}_{1\text{-}4}$ alkoxy, $\text{C}_{1\text{-}4}$ haloalkyl, $\text{C}_{3\text{-}6}$ cycloalkyl,

 C_{1-4} alkyl substituted with 0-1 R^{43} ,

aryl substituted with 0-3 R^{44} , and

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 $\rm R^{44}$:

 \mbox{R}^{43} is \mbox{C}_{3-6} cycloalkyl or aryl substituted with 0-3 $\mbox{R}^{44};$

 R^{44} , at each occurrence, is independently selected from H, halo, -OH, $NR^{46}R^{47}$, CO_2H , SO_2R^{45} , -CF₃, -OCF₃, -CN, - NO_2 , C_{1-4} alkyl, and C_{1-4} alkoxy;

30 R45 is C1-4 alkyl;

 R^{46} , at each occurrence, is independently selected from H and C_{1-4} alkyl;

35 R^{47} , at each occurrence, is independently selected from H, C_{1-4} alkyl, $-C(=0)NH(C_{1-4}$ alkyl), $-SO_2(C_{1-4}$ alkyl),

```
-SO_2 (phenyl), -C (=0) O (C_{1-4} alkyl), -C (=0) ( C_{1-4} alkyl),
            and -C(=0)H:
     R^{48}, at each occurrence, is independently selected from H,
  5
           C_{1-4} alkyl, -C(=0)NH(C_{1-4} alkyl), -C(=0)O(C_{1-4} alkyl),
            -C(=0)(C_{1-4} \text{ alkyl}), \text{ and } -C(=0)H;
     n is 1 or 2;
     m is 1 or 2; and
10
     n plus m is 2, 3, or 4;
     provided when n is 1, m is 2, and \mathbb{R}^7, \mathbb{R}^8, and \mathbb{R}^9 are
     independently selected from H, halogen, C_{1-4} alkyl, C_{1-4}
     alkoxy, C_{1-4} alkylthio or trifluoromethyl; then X is not a
15
     bond.
     14. A compound of Claim 12 wherein:
     \text{X is -CH}_2-, -O-, -S-, -CH_2CH_2-, -OCH_2-, -SCH_2-, -CH_2O-,
           or -CH2S-:
     R1 is selected from
            C_{2-5} alkyl substituted with Z,
            C_{2-5} alkenyl substituted with Z,
25
            C_{2-5} alkynyl substituted with Z,
            C3-6 cycloalkyl substituted with Z,
            aryl substituted with Z,
            5-6 membered heterocyclic ring system containing at
                least one heteroatom selected from the group
30
                consisting of N, O, and S, said heterocyclic ring
                system substituted with Z:
            C_{1-5} alkyl substituted with 0-2 R^2,
            C_{2-5} alkenyl substituted with 0-2 \mathbb{R}^2, and
            C2-5 alkynyl substituted with 0-2 R2;
35
     Z is selected from H.
           -CH (OH) R2,
```

```
th
O
Barry Maril Life
1.7
111
The sale of the
1 12
```

```
-C(ethylenedioxy)R2,
            -OR2.
            -SR2.
            -NR^2R^3,
  5
            -C(0)R^{2}.
            -C(O)NR2R3.
            -NR^3C(0)R^2,
            -C(O)OR2,
            -OC(O)R2.
10
            -CH (=NR4) NR2R3.
            -NHC(=NR^4)NR^2R^3,
            -S(0)R2.
            -S(0) 2R2,
            -S(0)_2NR^2R^3, and -NR^3S(0)_2R^2;
15
      {\ensuremath{\mathsf{R}}}^2, at each occurrence, is independently selected from
            C_{1-4} alkyl,
            C2-4 alkenyl,
            C2-4 alkynyl,
            C3-6 cycloalkyl.
            aryl substituted with 0-5 R42;
            C_{3-10} carbocyclic residue substituted with 0-3 R^{41}, and
            5-10 membered heterocyclic ring system containing from
                  1-4 heteroatoms selected from the group
25
                  consisting of N, O, and S substituted with 0-3
                  R41;
     \ensuremath{\mathbb{R}}^3, at each occurrence, is independently selected from
             H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, and
30
           C1-4 alkoxv:
     alternatively, \mathbb{R}^2 and \mathbb{R}^3 join to form a 5- or 6-membered
           ring optionally substituted with -O- or -N(R^4)-;
     {\ensuremath{\mathsf{R}}}^4, at each occurrence, is independently selected from H,
```

methyl, ethyl, propyl, and butyl;

```
R6a is H or C1-4 alkyl;
       R6b is H;
      alternatively, R^{6a} and R^{6b} are taken together to form =0 or
             =S:
      \mathbb{R}^7, \mathbb{R}^8, and \mathbb{R}^9, at each occurrence, are independently
              selected from
 10
             H, halo, -CF3, -OCF3, -OH, -OCH3, -CN, -NO2, -NR46R47,
             C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-4} haloalkyl,
                    C<sub>1-6</sub> alkoxy, (C<sub>1-4</sub> haloalkyl)oxy,
             C_{1-4} alkyl substituted with 0-2 R^{11}.
             C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
15
             aryl substituted with 0-5 R33.
             5-10 membered heterocyclic ring system containing from
                    1-4 heteroatoms selected from the group
                    consisting of N, O, and S substituted with 0-3
                    R31;
20
             OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13}.
             NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, CH(=NR^{14})NR^{12}R^{13}.
             {\rm NHC}\,(={\rm NR}^{14})\,{\rm NR}^{12}{\rm R}^{13},\ {\rm S}\,({\rm O})\,{\rm R}^{12},\ {\rm S}\,({\rm O})\,{}_2{\rm R}^{12},\ {\rm S}\,({\rm O})\,{}_2{\rm NR}^{12}{\rm R}^{13},
             NR^{14}S(O)_2R^{12}, NR^{14}S(O)_R^{12}, NR^{14}S(O)_2R^{12}, NR^{12}C(O)_R^{15},
25
             NR^{12}C(0)OR^{15}, NR^{12}S(0)_2R^{15}, and NR^{12}C(0)NHR^{15};
      R11 is selected from
            H, halo, -CF_3, -OCF_3, -OH, -OCH_3, -CN, -NO_2, -NR^{46}R^{47},
            C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-4} haloalkyl,
30
                   C_{1-6} alkoxy, (C_{1-4} haloalkyl)oxy,
            C_{3-10} carbocyclic residue substituted with 0-3 R^{33},
             aryl substituted with 0-5 R33,
            5-10 membered heterocyclic ring system containing from
                   1-4 heteroatoms selected from the group
35
                   consisting of N, O, and S substituted with 0-3 \,
```

R31.

10

15

```
OR^{12}, SR^{12}, NR^{12}R^{13}, C(O)H, C(O)R^{12}, C(O)NR^{12}R^{13},
               NR^{14}C(0)R^{12}, C(0)OR^{12}, OC(0)R^{12}, CH(=NR^{14})NR^{12}R^{13}
               NHC(=NR^{14})NR^{12}R^{13}, S(O)R^{12}, S(O)_2R^{12}, S(O)_2NR^{12}R^{13},
               and NR^{14}S(0)_{2}R^{12};
\mathbb{R}^{12}, at each occurrence, is independently selected from
```

 C_{1-4} alkyl, C2-4 alkenyl,

C2-4 alkynyl,

C3-6 cycloalkyl,

phenyl substituted with 0-5 R33;

 C_{3-10} carbocyclic residue substituted with 0-3 \mathbb{R}^{33} , and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3

R31.

- \mathbb{R}^{13} , at each occurrence, is independently selected from H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;
- alternatively, ${\ensuremath{\mathsf{R}}}^{12}$ and ${\ensuremath{\mathsf{R}}}^{13}$ join to form a 5- or 6-membered ring optionally substituted with -O- or $-N(R^{14})$ -;
- \mathbb{R}^{14} , at each occurrence, is independently selected from H 25 and C_{1-4} alkyl;
 - \mathbb{R}^{31} , at each occurrence, is independently selected from H, OH, halo, CF3, methyl, and ethyl;
- \mathbb{R}^{33} , at each occurrence, is independently selected from 30 H, OH, halo, CN, NO2, CF3, methyl, and ethyl;
- \mathbf{R}^{41} , at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO₂R⁴⁵, NR⁴⁶R⁴⁷, NO₂, CN, =0, 35 C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl, C1-4 alkyl substituted with 0-1 R43, aryl substituted with 0-3 R42, and

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 $\rm R^{44}$;

5

10

15

25

35

 R^{42} , at each occurrence, is independently selected from H, CF₃, halo, OH, CO_2H , SO_2R^{45} , SR^{45} , $NR^{46}R^{47}$, OR^{48} , NO_2 , CN, CH(=NH) NH_2 , NHC(=NH) NH_2 ,

C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₄ alkoxy, C₁₋₄ haloalkyl, C₃₋₆ cycloalkyl.

C1-4 alkyl substituted with 0-1 R43.

aryl substituted with 0-3 R44, and

5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3 \mathbb{R}^{44} :

 R^{43} is C_{3-6} cycloalkyl or aryl substituted with 0-3 R^{44} ;

20 R^{44} , at each occurrence, is independently selected from H, halo, -OH, $NR^{46}R^{47}$, CO_2H , SO_2R^{45} , -CF₃, -OCF₃, -CN, - NO_2 , C_{1-4} alkyl, and C_{1-4} alkoxy;

 \mathbb{R}^{45} is \mathbb{C}_{1-4} alkyl;

 $R^{46}, \mbox{ at each occurrence, is independently selected from H } \mbox{ and } C_{1-3} \mbox{ alkyl;}$

 R^{47} , at each occurrence, is independently selected from H, 30 C_{1-4} alkyl, $-C(=0)NH(C_{1-4}$ alkyl), $-SO_2(C_{1-4}$ alkyl), $-SO_2(phenyl)$, $-C(=0)O(C_{1-4}$ alkyl), $-C(=0)(C_{1-4}$ alkyl), and -C(=0)H:

 $R^{48},$ at each occurrence, is independently selected from H, $C_{1-4} \text{ alkyl}, \text{ -C(=0)NH(C}_{1-4} \text{ alkyl}), \text{ -C(=0)O(C}_{1-4} \text{ alkyl}), \\ \text{-C(=0)(} C_{1-4} \text{ alkyl}), \text{ and -C(=0)H;}$

```
m is 1 or 2: and
         n plus m is 2, 3, or 4.
        15. A compound of Claim 13 wherein:
        X is -CH2-, -O- or -S-;
        R1 is selected from
   10
               C2-4 alkyl substituted with Z,
               C_{2-4} alkenyl substituted with Z,
               C2-4 alkynyl substituted with Z.
               C<sub>3-6</sub> cycloalkyl substituted with Z,
               aryl substituted with Z,
               5-6 membered heterocyclic ring system containing at
(1 15
                    least one heteroatom selected from the group
                    consisting of N, O, and S, said heterocyclic ring
                    system substituted with Z;
               C_{2-4} alkyl substituted with 0-2 \mathbb{R}^2, and
  20
               C2-4 alkenyl substituted with 0-2 R2;
        Z is selected from H.
              -CH (OH) R2,
              -C(ethylenedioxy)R2,
   25
              -OR2.
              -SR2.
              -NR^2R^3.
              -C(O)R2,
              -C(O)NR2R3,
              -NR^3C(O)R^2.
   30
              -C(O)OR2,
              -S(0)R2,
              -S(0)_2R^2,
              -S(0)_2NR^2R^3, and -NR^3S(0)_2R^2;
   35
        \ensuremath{\mathbb{R}}^2, at each occurrence, is independently selected from
             phenyl substituted with 0-5 R42;
```

n is 1 or 2;

l. ... (1)

13

13

11

1.7 0 \$.i

```
C_{3-10} carbocyclic residue substituted with 0-3 R^{41}, and
           5-10 membered heterocyclic ring system containing from
                 1-4 heteroatoms selected from the group
                 consisting of N, O, and S substituted with 0-3
 5
                 R41;
     \mathbb{R}^3, at each occurrence, is independently selected from
            H, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, and
           C_{1-4} alkoxy;
10
     alternatively, \mathbb{R}^2 and \mathbb{R}^3 join to form a 5- or 6-membered
           ring optionally substituted with -O- or -N(R4)-;
     \mathbb{R}^4, at each occurrence, is independently selected from H,
15
           methyl, ethyl, propyl, and butyl;
     R6a is H or C1-4 alkyl;
     R6b is H;
20
     alternatively, R^{6a} and R^{6b} are taken together to form =0 or
           =S;
     R^7, R^8, and R^9, at each occurrence, are independently
25
           selected from
           H, halo, -CF3, -OCF3, -OH, -OCH3, -CN, -NO2,
           C_{1-4} alkyl, C_{1-4} haloalkyl, C_{1-4} alkoxy, (C_{1-3})
                 haloalkyl)oxy, and
           C_{1-4} alkyl substituted with 0-2 R^{11};
30
     R11 is selected from
           H, halo, -CF3, -OCF3, -OH, -OCH3, -CN, -NO2,
           C_{1-4} alkyl, C_{1-4} haloalkyl, C_{1-4} alkoxy, and (C_{1-3})
                haloalkyl)oxy;
35
```

 \mathbb{R}^{33} , at each occurrence, is independently selected from H, OH, halo, \mathbb{CF}_3 , and methyl;

10

 \mathbb{R}^{41} , at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO₂R⁴⁵, NR⁴⁶R⁴⁷, NO₂, CN, =0. C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl, C1-4 alkyl substituted with 0-1 R43, aryl substituted with 0-3 R42, and 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3

R42, at each occurrence, is independently selected from H, CF₃, halo, OH, CO₂H, SO₂R⁴⁵, SR⁴⁵, NR⁴⁶R⁴⁷, OR⁴⁸, NO2, CN, CH(=NH)NH2, NHC(=NH)NH2,

 C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-4} alkoxy, C_{1-4} haloalkyl, C3-6 cycloalkyl,

C₁₋₄ alkyl substituted with 0-1 R⁴³,

aryl substituted with $0-3\ R^{44}$, and

- 5-10 membered heterocyclic ring system containing from 1-4 heteroatoms selected from the group consisting of N, O, and S substituted with 0-3R44;
- R43 is cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, 25 phenyl, or pyridyl, each substituted with 0-3 R44;
 - \mathbb{R}^{44} , at each occurrence, is independently selected from H, halo, -OH, $NR^{46}R^{47}$, CO_2H , SO_2R^{45} , -CF₃, -OCF₃, -CN, -NO2, methyl, ethyl, propyl, butyl, methoxy, ethoxy, propoxy, and butoxy;
- 30

R44:

- R45 is methyl, ethyl, propyl, or butyl;
- ${\rm R}^{46}$, at each occurrence, is independently selected from H, 35 methyl, ethyl, propyl, and butyl:
 - ${\ensuremath{R^{47}}}$, at each occurrence, is independently selected from

```
- It giest just gemein gemein gemein zwenn im gemein ist gemein is
```

```
H, methyl, ethyl, n-propyl, i-propyl, n-butyl,
           i-butyl, -C(=0)NH(methyl), -C(=0)NH(ethyl),
           -SO_2 (methyl), -SO_2 (ethyl), -SO_2 (phenyl),
           -C(=0)O(methyl), -C(=0)O(ethyl), -C(=0)(methyl),
           -C(=0) (ethyl), and -C(=0)H;
 5
     R^{48}, \ \text{at each occurrence, is independently selected from }
          H, methyl, ethyl, n-propyl, i-propyl, -
          C(=0) NH (methyl), -C(=0) NH (ethyl), -C(=0) O (methyl), -
10
          C(=0)O(ethyl), -C(=0)(methyl), -C(=0)(ethyl), and -
          C(=0)H;
     n is 1 or 2;
     m is 1 or 2; and
     n plus m is 2 or 3.
     16. A compound of Claim 13 wherein:
     X is -CH2-, -O- or -S-;
20
     R1 is selected from
          ethyl substituted with Z,
          propyl substituted with Z.
          butyl substituted with Z,
25
          propenyl substituted with Z,
          butenyl substituted with Z,
          ethyl substituted with R2,
          propyl substituted with R2.
          butyl substituted with R2.
30
          propenyl substituted with R2, and
          butenyl substituted with R2;
     Z is selected from H,
          -CH (OH) R2.
35
          -OR2.
          -SR2.
          -NR2R3.
```

```
the first party and party and the same party and the same of the same of the same and the same a
```

```
-C(0)R2.
           -C(0)NR2R3.
           -NR^3C(0)R^2
           -C(0)OR2.
 5
           -S(0)R2,
           -S(0)2R2,
           -S(0)_2NR^2R^3, and -NR^3S(0)_2R^2;
     {\ensuremath{\mathsf{R}}}^2, at each occurrence, is independently selected from
10
           phenyl substituted with 0-3 R42;
           naphthyl substituted with 0-3 R42;
           cyclopropyl substituted with 0-3 R41;
           cyclobutyl substituted with 0-3 R41;
           cyclopentyl substituted with 0-3 R41;
15
           cyclohexyl substituted with 0-3 R41;
          pyridyl substituted with 0-3 R41;
           indolyl substituted with 0-3 R41:
           indolinyl substituted with 0-3 R41;
          benzimidazolyl substituted with 0-3 R41;
20
          benzotriazolyl substituted with 0-3 R41:
          benzothienyl substituted with 0-3 R41;
          benzofuranyl substituted with 0-3 R41;
          phthalimid-1-yl substituted with 0-3 R41;
          inden-2-yl substituted with 0-3 R41;
          2,3-dihydro-1H-inden-2-yl substituted with 0-3 R41;
25
          indazolyl substituted with 0-3 R41;
          tetrahydroquinolinyl substituted with 0-3 R41; and
          tetrahydro-isoquinolinyl substituted with 0-3 R41;
30
    {\ensuremath{\mathsf{R}}}^3, at each occurrence, is independently selected from
           H, methyl, and ethyl;
    R6a is H or C1-4 alkyl;
35
   R6b is H:
```

25

- alternatively, R6a and R6b are taken together to form =0 or =S:
- R^7 , R^8 , and R^9 , at each occurrence, are independently 5 selected from H, F, Cl, methyl, ethyl, methoxy, -CF3, and -OCF3;
- ${\bf R}^{41}$, at each occurrence, is independently selected from H, F, Cl, Br, OH, CF3, NO2, CN, =0, methyl, ethyl, 10 propyl, butyl, methoxy, and ethoxy;
 - \mathbb{R}^{42} , at each occurrence, is independently selected from H, F, Cl, Br, OH, CF₃, SO₂R⁴⁵, SR⁴⁵, NR⁴⁶R⁴⁷, OR⁴⁸, NO₂, CN, =0, methyl, ethyl, propyl, butyl, methoxy, and ethoxy;
 - R^{45} is methyl, ethyl, propyl, or butyl;
 - ${\bf R}^{46}$, at each occurrence, is independently selected from H, methyl, ethyl, propyl, and butyl;
 - \mathbb{R}^{47} , at each occurrence, is independently selected from H, methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, -C(=O)NH(methyl), -C(=O)NH(ethyl), $-SO_2$ (methyl), $-SO_2$ (ethyl), $-SO_2$ (phenyl), $-C(=0) \circ (methyl), -C(=0) \circ (ethyl), -C(=0) (methyl),$ -C(=0) (ethyl), and -C(=0)H:
- ${\bf R}^{48},$ at each occurrence, is independently selected from 30 H, methyl, ethyl, n-propyl, i-propyl, -C(=0) NH (methyl), -C(=0) NH (ethyl), -C(=0) O (methyl), -C(=0)C(=0)O(ethyl), -C(=0)(methyl), -C(=0)(ethyl), and -C(=0)H:
- 35 n is 1; and m is 1.

```
17. A compound of Claim 13 of Formula (II)
```

wherein:

5

PI

6/1

The last

```
b is a single bond wherein the bridging hydrogens are
either cis or trans;
```

```
R1 is selected from
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-fluoro-phenyl),
10
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-bromo-phenv1).
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-methyl-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-methoxy-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-(3,4-dichloro-phenyl)phenyl),
15
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(3-methyl-4-fluoro-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2,3-dimethoxy-phenyl).
            -(CH<sub>2</sub>)<sub>3</sub>C(=O)(phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-chloro-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=O)(3-methyl-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-t-butyl-phenyl),
20
           -(CH<sub>2</sub>)<sub>3</sub>C(=O)(3,4-difluoro-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-methoxy-5-fluoro-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-fluoro-1-naphthyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(benzyl),
25
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(4-pyridyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=O)(3-pyridy1),
           -(CH2)3CH(OH)(4-fluoro-phenyl),
           -(CH2)3CH(OH)(4-pyridyl),
           -(CH2)3CH(OH)(2,3-dimethoxy-phenyl),
3.0
           -(CH2)3S(3-fluoro-phenyl),
           -(CH2)3S(4-fluoro-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>S(=0)(4-fluoro-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>SO<sub>2</sub>(3-fluoro-phenyl),
           -(CH2)3SO2(4-fluoro-phenyl),
                                                   -317-
```

```
-(CH2)30(4-fluoro-phenyl),
                 - (CH2) 30 (phenyl),
                 -(CH2)30(3-pyridyl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(4-pyridyl),
       5
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-phenyl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-5-F-phenvl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-4-F-phenv1).
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-3-F-phenyl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-4-Cl-phenvl).
     10
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-4-OH-phenyl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NH<sub>2</sub>-4-Br-phenyl),
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NHC(=O)Me-4-F-phenyl),
Carl Start Start Start
                 -(CH<sub>2</sub>)<sub>3</sub>O(2-NHC(=O)Me-phenyl),
                 -(CH2)3NH(4-fluoro-phenyl),
                 -(CH2)3N(methyl)(4-fluoro-phenyl),
     15
0
                 -(CH2)3CO2(ethv1).
                 -(CH<sub>2</sub>)<sub>3</sub>C(=O)N(methyl)(methoxy).
11
                 -(CH<sub>2</sub>)<sub>3</sub>C(=0)NH(4-fluoro-phenyl).
1.1
                 -(CH<sub>2</sub>)<sub>2</sub>NHC(=O)(phenyl),
111
1 1
    20
                 -(CH<sub>2</sub>)<sub>2</sub>NMeC(=O)(phenv1).
0
                 -(CH<sub>2</sub>)<sub>2</sub>NHC(=O)(2-fluoro-phenvl).
                 -(CH2)2NMeC(=0)(2-fluoro-phenyl),
                 -(CH2)2NHC(=0)(4-fluoro-phenyl),
                -(CH2)2NMeC(=O)(4-fluoro-phenyl),
    25
                -(CH2)2NHC(=0)(2,4-difluoro-phenyl).
                -(CH2)2NMeC(=0)(2,4-difluoro-phenyl).
                -(CH<sub>2</sub>)<sub>3</sub>(3-indoly1),
                -(CH<sub>2</sub>)<sub>3</sub>(1-methyl-3-indolvl).
                -(CH<sub>2</sub>)<sub>3</sub>(1-indoly1),
                -(CH<sub>2</sub>)<sub>3</sub>(1-indolinyl),
    30
                -(CH<sub>2</sub>)<sub>3</sub>(1-benzimidazolv1).
                -(CH<sub>2</sub>)<sub>3</sub>(1H-1,2,3-benzotriazol-1-yl),
                -(CH<sub>2</sub>)<sub>3</sub>(1H-1,2,3-benzotriazol-2-yl),
                -(CH<sub>2</sub>)<sub>2</sub>(1H-1,2,3-benzotriazol-1-yl),
    35
                -(CH<sub>2</sub>)<sub>2</sub>(1H-1,2,3-benzotriazol-2-vl).
                -(CH2)3(3,4 dihydro-1(2H)-quinolinyl),
                -(CH<sub>2</sub>)<sub>2</sub>C(=0)(4-fluoro-phenyl),
```

```
-(CH<sub>2</sub>)<sub>2</sub>C(=O)NH(4-fluoro-phenyl),
            -CH2CH2 (3-indoly1),
            -CH2CH2 (1-phthalimidyl),
            -(CH2)4C(=0)N(methyl)(methoxy),
  5
            -(CH<sub>2</sub>)<sub>4</sub>CO<sub>2</sub>(ethyl),
            -(CH<sub>2</sub>)<sub>4</sub>C(=0)(phenyl),
            - (CH2)4 (cvclohexvl).
            -(CH2)3CH(pheny1)2,
            -CH2CH2CH=C(phenyl)2.
            -CH2CH2CH=CMe(4-F-phenyl),
 10
            -(CH2)3CH(4-fluoro-phenyl)2.
            -CH2CH2CH=C(4-fluoro-phenyl)2,
            -(CH<sub>2</sub>)<sub>2</sub>(2,3-dihydro-1H-inden-2-v1).
            -(CH_2)_3C(=0)(2-NH_2-phenyl),
15
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-NH<sub>2</sub>-5-F-phenv1).
            -(CH_2)_3C(=0)(2-NH_2-4-F-phenyl),
            -(CH_2)_3C(=0)(2-NH_2-3-F-pheny1),
            -(CH_2)_3C(=0)(2-NH_2-4-Cl-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=O)(2-NH<sub>2</sub>-4-OH-phenyl),
20
            -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-NH<sub>2</sub>-4-Br-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>(1H-indazol-3-yl),
            -(CH<sub>2</sub>)<sub>3</sub>(5-F-1H-indazol-3-v1).
            -(CH<sub>2</sub>)<sub>3</sub>(7-F-1H-indazol-3-yl),
            -(CH<sub>2</sub>)<sub>3</sub>(6-Cl-1H-indazol-3-yl),
25
            -(CH<sub>2</sub>)<sub>3</sub>(6-Br-1H-indazol-3-yl),
            -(CH<sub>2</sub>)<sub>3</sub>C(=O)(2-NHMe-phenyl),
            -(CH<sub>2</sub>)<sub>3</sub>(1-benzothien-3-v1).
            -(CH<sub>2</sub>)<sub>3</sub>(6-F-1H-indol-1-vl),
            -(CH<sub>2</sub>)<sub>3</sub>(5-F-1H-indol-1-yl),
30
           -(CH<sub>2</sub>)<sub>3</sub>(6-F-2,3-dihydro-1H-indol-1-vl).
           -(CH<sub>2</sub>)<sub>3</sub>(5-F-2,3-dihydro-1H-indol-1-yl),
           -(CH<sub>2</sub>)<sub>3</sub>(6-F-1H-indol-3-v1),
           -(CH<sub>2</sub>)<sub>3</sub>(5-F-1H-indol-3-v1).
           -(CH<sub>2</sub>)<sub>3</sub>(5-F-1H-indol-3-y1),
35
           -(CH<sub>2</sub>)<sub>3</sub>(9H-purin-9-v1).
           -(CH2)3(7H-purin-7-v1).
           -(CH<sub>2</sub>)<sub>3</sub>(6-F-1H-indazol-3-v1).
```

(3

219

ΓÜ

in

4.3

7

-0 111

16

13

```
-(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-NHSO<sub>2</sub>Me-4-F-phenyl),
           -(CH_2)_3C(=0)(2-NHC(=0)Me-4-F-pheny1),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-NHC(=0)Me-phenyl),
           -(CH_2)_3C(=0)(2-NHCO_2Et-4-F-phenyl),
 5
           -(CH_2)_3C(=0)(2-NHC(=0)NHEt-4-F-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-NHCHO-4-F-phenv1).
          -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-OH-4-F-phenyl),
           -(CH<sub>2</sub>)<sub>3</sub>C(=0)(2-MeS-4-F-phenyl),
          -(CH_2)_3C(=0)(2-NHSO_2Me-4-F-phenyl),
10
          - (CH2) 2C (Me) CO2Me,
          -(CH<sub>2</sub>)<sub>2</sub>C(Me)CH(OH)(4-F-phenyl)<sub>2</sub>
          -(CH<sub>2</sub>)<sub>2</sub>C(Me)CH(OH)(4-Cl-phenyl)<sub>2</sub>,
          -(CH_2)_2C(Me)C(=0)(4-F-phenyl),
          -(CH_2)_2C(Me)C(=0)(2-MeO-4-F-phenyl),
15
          -(CH<sub>2</sub>)<sub>2</sub>C(Me)C(=0)(3-Me-4-F-phenv1).
          -(CH_2)_2C(Me)C(=0)(2-Me-phenyl),
          -(CH<sub>2</sub>)<sub>2</sub>C(Me)C(=0)phenyl,
20
                                                                                     and
```

 $25~~R^{7},~R^{8},~{\rm and}~R^{9},~{\rm at}~{\rm each}~{\rm occurrence},~{\rm are}~{\rm independently}$ selected from

hydrogen, fluoro, chloro, bromo, cyano, methyl, ethyl, propyl, isopropyl, butyl, t-butyl, nitro, trifluoromethyl, methoxy, ethoxy, isopropoxy,

30 trifluoromethoxy, phenyl, benzyl,

```
HC(=0) -, methylC(=0) -, ethylC(=0) -, propylC(=0) -, isopropylC(=0) -, n-butylC(=0) -, isobutylC(=0) -, secbutylC(=0) -, tertbutylC(=0) -, phenylC(=0) -,
```

5

methylC(=0)NH-, ethylC(=0)NH -, propylC(=0)NH-,
isopropylC(=0)NH-, n-butylC(=0)NH-, isobutylC(=0)NH-,
secbutylC(=0)NH-, tertbutylC(=0)NH-, phenylC(=0)NH-,

methylamino-, ethylamino-, propylamino-, isopropylamino-,
n-butylamino-, isobutylamino-, secbutylamino-,
tertbutylamino-, phenylamino-,

provided that two of substituents R^7 , R^8 , and R^9 , are independently selected from hydrogen, fluoro, chloro, bromo, cyano, methyl, ethyl, propyl, isopropyl, butyl, t-butyl, nitro, trifluoromethyl, methoxy, ethoxy, isopropoxy, and trifluoromethoxy.

- 20 18. A compound selected from the group consisting of compounds disclosed in Table 1.
 - 19. A compound selected from the group consisting of compounds disclosed in Table 2.

25

- 20. A compound selected from the group consisting of compounds disclosed in Table 3.
- 21. A pharmaceutical composition comprising a 30 pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.
- 22. A method for treating a human suffering from a 35 disorder associated with 5HT2C receptor modulation comprising administering to a patient in need thereof a

therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.

- 23. A method of Claim 22 for treating a human suffering from a disorder associated with 5HT2C receptor modulation wherein the compound is a 5HT2C agonist.
- 24. A method for treating a human suffering from a disorder associated with 5HT2A receptor modulation comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.
- 25. A method of Claim 24 for treating a human suffering 5 from a disorder associated with 5HT2A receptor modulation wherein the compound is a 5HT2A antagonist.
 - 26. A method for treating obesity comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.
 - 27. A method for treating schizophrenia comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.
- 28. A method for treating depression comprising administering to a patient in need thereof a 30 therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.